

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
 Data File : BF127171.D
 Acq On : 03 Mar 2022 15:05
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 03/04/2022
 Supervised By : Jagrut Upadhyay 03/04/2022

Quant Time: Mar 03 17:01:46 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Mar 01 23:57:09 2022

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
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Internal Standards

1) 1,4-Dichlorobenzene-d4	6.893	152	104344	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	415007	20.000	ng	# 0.00
39) Acenaphthene-d10	9.934	164	212567	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	386648	20.000	ng	0.00
76) Chrysene-d12	14.074	240	256942	20.000	ng	# 0.00
86) Perylene-d12	15.563	264	190366	20.000	ng	0.00

03/04/2022

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System Monitoring Compounds

5) 2-Fluorophenol	5.510	112	547238	81.058	ng	0.00
7) Phenol-d6	6.522	99	734229	80.599	ng	0.00
23) Nitrobenzene-d5	7.457	82	701459	76.293	ng	0.00
42) 2,4,6-Tribromophenol	10.728	330	206177	84.243	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	1127032	78.054	ng	0.00
79) Terphenyl-d14	13.010	244	1294747	74.076	ng	0.00

Target Compounds

Compound	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.658	88	106418	34.448	ng	98
3) Pyridine	3.428	79	329367	40.648	ng	97
4) n-Nitrosodimethylamine	3.399	42	148402	37.387	ng	83
6) Aniline	6.557	93	426032	39.471	ng	97
8) 2-Chlorophenol	6.675	128	296180	40.888	ng	97
9) Benzaldehyde	6.446	77	198043	35.714	ng	99
10) Phenol	6.534	94	399614	43.414	ng	86
11) bis(2-Chloroethyl)ether	6.628	93	284090	39.350	ng	98
12) 1,3-Dichlorobenzene	6.834	146	320658	41.037	ng	98
13) 1,4-Dichlorobenzene	6.910	146	321441	40.577	ng	98
14) 1,2-Dichlorobenzene	7.063	146	301962	39.972	ng	98
15) Benzyl Alcohol	7.034	79	273153	35.592	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.163	45	386301	33.980	ng	90
17) 2-Methylphenol	7.140	107	254670	40.636	ng	94
18) Hexachloroethane	7.398	117	120985	39.844	ng	95
19) n-Nitroso-di-n-propyla...	7.304	70	214023	36.460	ng	96
20) 3+4-Methylphenols	7.298	107	319036	39.203	ng	91
22) Acetophenone	7.304	105	408194	37.357	ng	# 95
24) Nitrobenzene	7.475	77	334925	38.561	ng	95
25) Isophorone	7.710	82	566687	37.247	ng	99
26) 2-Nitrophenol	7.792	139	154733	41.842	ng	# 90
27) 2,4-Dimethylphenol	7.822	122	217122	35.739	ng	91
28) bis(2-Chloroethoxy)met...	7.922	93	355384	38.530	ng	99
29) 2,4-Dichlorophenol	8.034	162	231870	40.596	ng	99
30) 1,2,4-Trichlorobenzene	8.116	180	253140	39.642	ng	98
31) Naphthalene	8.198	128	846823	39.089	ng	100
32) Benzoic acid	7.945	122	155479m	36.087	ng	
33) 4-Chloroaniline	8.245	127	340107	39.885	ng	97
34) Hexachlorobutadiene	8.304	225	145416	39.007	ng	98
35) Caprolactam	8.628	113	81486	40.854	ng	# 80
36) 4-Chloro-3-methylphenol	8.728	107	288438	39.514	ng	98
37) 2-Methylnaphthalene	8.887	142	544981	38.768	ng	98
38) 1-Methylnaphthalene	8.987	142	532119	39.529	ng	99
40) 1,2,4,5-Tetrachloroben...	9.051	216	228963	40.919	ng	97
41) Hexachlorocyclopentadiene	9.034	237	108159	35.640	ng	94
43) 2,4,6-Trichlorophenol	9.169	196	164524	40.089	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.210	196	174988	40.519	ng	94	03/04/2022
46) 1,1'-Biphenyl	9.351	154	658381	39.093	ng	98	Supervised By : Jagrut Upadhyay
47) 2-Chloronaphthalene	9.381	162	491704	39.371	ng	95	
48) 2-Nitroaniline	9.481	65	193907	39.346	ng	94	
49) Acenaphthylene	9.798	152	799067	39.389	ng	98	
50) Dimethylphthalate	9.657	163	597307	39.312	ng	99	
51) 2,6-Dinitrotoluene	9.722	165	128459	41.555	ng	99	03/04/2022
52) Acenaphthene	9.969	154	519452m	39.372	ng		
53) 3-Nitroaniline	9.892	138	155169	41.064	ng	94	
54) 2,4-Dinitrophenol	9.998	184	65065	39.766	ng	85	
55) Dibenzofuran	10.139	168	716529	39.496	ng	92	
56) 4-Nitrophenol	10.051	139	107605	38.556	ng	# 86	
57) 2,4-Dinitrotoluene	10.122	165	174268	41.694	ng	# 98	
58) Fluorene	10.486	166	559268	39.577	ng	100	
59) 2,3,4,6-Tetrachlorophenol	10.257	232	138504	40.454	ng	92	
60) Diethylphthalate	10.351	149	606624	38.209	ng	99	
61) 4-Chlorophenyl-phenyle...	10.475	204	262001	39.319	ng	97	
62) 4-Nitroaniline	10.510	138	158327	42.431	ng	# 82	
63) Azobenzene	10.633	77	665833	37.663	ng	96	
65) 4,6-Dinitro-2-methylph...	10.533	198	96989	42.692	ng	98	
66) n-Nitrosodiphenylamine	10.592	169	485409	38.286	ng	98	
67) 4-Bromophenyl-phenylether	10.963	248	168451	38.993	ng	91	
68) Hexachlorobenzene	11.028	284	184946	39.768	ng	93	
69) Atrazine	11.122	200	153481	39.031	ng	95	
70) Pentachlorophenol	11.228	266	88123	34.712	ng	96	
71) Phenanthrene	11.451	178	842198	38.916	ng	99	
72) Anthracene	11.504	178	836052	38.806	ng	99	
73) Carbazole	11.657	167	783064	39.634	ng	99	
74) Di-n-butylphthalate	11.980	149	940840	37.042	ng	100	
75) Fluoranthene	12.639	202	835338	40.652	ng	97	
77) Benzidine	12.763	184	244988	35.086	ng	97	
78) Pyrene	12.869	202	843102	37.769	ng	99	
80) Butylbenzylphthalate	13.480	149	386348	36.559	ng	92	
81) Benzo(a)anthracene	14.063	228	681188	39.325	ng	99	
82) 3,3'-Dichlorobenzidine	14.021	252	222071	40.680	ng	# 98	
83) Chrysene	14.098	228	637782	39.154	ng	99	
84) Bis(2-ethylhexyl)phtha...	14.033	149	482206	34.755	ng	96	
85) Di-n-octyl phthalate	14.651	149	688298	34.455	ng	95	
87) Indeno(1,2,3-cd)pyrene	17.074	276	500706	40.113	ng	# 88	
88) Benzo(b)fluoranthene	15.121	252	498588	38.996	ng	# 95	
89) Benzo(k)fluoranthene	15.151	252	517867	41.993	ng	# 96	
90) Benzo(a)pyrene	15.498	252	402653	40.318	ng	# 94	
91) Dibenzo(a,h)anthracene	17.092	278	409066	39.624	ng	# 93	
92) Benzo(g,h,i)perylene	17.539	276	416994	40.971	ng	# 89	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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