

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091924\
 Data File : BF139468.D
 Acq On : 19 Sep 2024 17:19
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/21/2024
 Supervised By :mohammad ahmed 09/22/2024

Quant Time: Sep 19 17:41:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 14 02:58:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	70449	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	249250	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	134416	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	256666	20.000	ng	0.00
76) Chrysene-d12	14.033	240	130668	20.000	ng	0.00
86) Perylene-d12	15.504	264	130863	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	323498	74.911	ng	0.00
7) Phenol-d6	6.510	99	422786	74.588	ng	0.00
23) Nitrobenzene-d5	7.445	82	396430	74.771	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	101218	70.731	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	693291	72.339	ng	0.00
79) Terphenyl-d14	12.980	244	666185	83.683	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.628	88	68852	39.740	ng	95
3) Pyridine	3.381	79	165072	43.213	ng	97
4) n-Nitrosodimethylamine	3.340	42	117209	35.400	ng	89
6) Aniline	6.540	93	183034	42.789	ng	96
8) 2-Chlorophenol	6.663	128	168423	37.888	ng	99
9) Benzaldehyde	6.428	77	118960	36.325	ng	98
10) Phenol	6.522	94	220996	38.281	ng	97
11) bis(2-Chloroethyl)ether	6.616	93	158187	36.051	ng	99
12) 1,3-Dichlorobenzene	6.822	146	187826	37.153	ng	98
13) 1,4-Dichlorobenzene	6.898	146	193083	37.794	ng	99
14) 1,2-Dichlorobenzene	7.051	146	180186	37.326	ng	99
15) Benzyl Alcohol	7.016	79	162414	35.914	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.151	45	271975	40.220	ng	100
17) 2-Methylphenol	7.128	107	134011	37.944	ng	97
18) Hexachloroethane	7.392	117	71079	36.528	ng	98
19) n-Nitroso-di-n-propyla...	7.287	70	122525	36.477	ng	100
20) 3+4-Methylphenols	7.281	107	175066	36.332	ng	97
22) Acetophenone	7.287	105	233969	35.906	ng	99
24) Nitrobenzene	7.463	77	206287	37.804	ng	98
25) Isophorone	7.698	82	314175	37.469	ng	99
26) 2-Nitrophenol	7.775	139	86369	40.626	ng	93
27) 2,4-Dimethylphenol	7.810	122	107137m	38.205	ng	
28) bis(2-Chloroethoxy)met...	7.910	93	185257	38.766	ng	99
29) 2,4-Dichlorophenol	8.016	162	134412	37.729	ng	99
30) 1,2,4-Trichlorobenzene	8.104	180	155644	37.496	ng	99
31) Naphthalene	8.181	128	496623	38.166	ng	100
32) Benzoic acid	7.916	122	100119m	38.150	ng	
33) 4-Chloroaniline	8.234	127	150085	37.702	ng	97
34) Hexachlorobutadiene	8.298	225	100248	35.548	ng	99
35) Caprolactam	8.598	113	44162	40.298	ng	95
36) 4-Chloro-3-methylphenol	8.710	107	151197	37.136	ng	96
37) 2-Methylnaphthalene	8.875	142	310342	37.575	ng	99
38) 1-Methylnaphthalene	8.975	142	301437	37.174	ng	99
40) 1,2,4,5-Tetrachloroben...	9.039	216	150411	37.757	ng	99
41) Hexachlorocyclopentadiene	9.022	237	48455	34.836	ng	98
43) 2,4,6-Trichlorophenol	9.151	196	97768	37.723	ng	99

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44) 2,4,5-Trichlorophenol	9.186	196	105325	38.526	ng	99
46) 1,1'-Biphenyl	9.333	154	391497	37.859	ng	99
47) 2-Chloronaphthalene	9.363	162	292951	37.888	ng	99
48) 2-Nitroaniline	9.457	65	110051	38.267	ng	99
49) Acenaphthylene	9.775	152	447342	38.435	ng	99
50) Dimethylphthalate	9.633	163	330365	37.447	ng	100
51) 2,6-Dinitrotoluene	9.698	165	76969	39.382	ng	95
52) Acenaphthene	9.945	154	302408	37.228	ng	99
53) 3-Nitroaniline	9.869	138	82265	42.118	ng	88
54) 2,4-Dinitrophenol	9.969	184	39809	36.554	ng	95
55) Dibenzofuran	10.122	168	427160	37.728	ng	96
56) 4-Nitrophenol	10.022	139	60541	37.459	ng	# 82
57) 2,4-Dinitrotoluene	10.098	165	104235	39.504	ng	97
58) Fluorene	10.463	166	346147	37.471	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.233	232	82892	36.653	ng	99
60) Diethylphthalate	10.333	149	333754	36.659	ng	98
61) 4-Chlorophenyl-phenyle...	10.451	204	163408	36.398	ng	96
62) 4-Nitroaniline	10.475	138	82531	38.716	ng	88
63) Azobenzene	10.610	77	334119	37.313	ng	97
65) 4,6-Dinitro-2-methylph...	10.504	198	56312	41.570	ng	95
66) n-Nitrosodiphenylamine	10.569	169	285219	38.009	ng	99
67) 4-Bromophenyl-phenylether	10.945	248	95229	37.440	ng	96
68) Hexachlorobenzene	11.004	284	109184	36.576	ng	99
69) Atrazine	11.092	200	61338	34.335	ng	97
70) Pentachlorophenol	11.198	266	66998	37.404	ng	97
71) Phenanthrene	11.422	178	469140	37.886	ng	99
72) Anthracene	11.474	178	462152	38.172	ng	99
73) Carbazole	11.627	167	448869	38.736	ng	99
74) Di-n-butylphthalate	11.957	149	484155	37.507	ng	100
75) Fluoranthene	12.610	202	508429	38.447	ng	99
77) Benzidine	12.733	184	104748	55.952	ng	99
78) Pyrene	12.839	202	514572	45.864	ng	100
80) Butylbenzylphthalate	13.451	149	157826	41.327	ng	99
81) Benzo(a)anthracene	14.021	228	324244	37.461	ng	100
82) 3,3'-Dichlorobenzidine	13.986	252	94960	37.463	ng	99
83) Chrysene	14.063	228	305611	38.379	ng	99
84) Bis(2-ethylhexyl)phtha...	14.010	149	181970	37.456	ng	99
85) Di-n-octyl phthalate	14.621	149	241988	33.822	ng	99
87) Indeno(1,2,3-cd)pyrene	16.974	276	332904	42.782	ng	97
88) Benzo(b)fluoranthene	15.074	252	292286	36.215	ng	99
89) Benzo(k)fluoranthene	15.104	252	265014	35.613	ng	99
90) Benzo(a)pyrene	15.439	252	250680	37.451	ng	99
91) Dibenzo(a,h)anthracene	16.992	278	268828	41.808	ng	98
92) Benzo(g,h,i)perylene	17.421	276	286406	44.986	ng	97

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

