

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072324\
 Data File : BF138629.D
 Acq On : 23 Jul 2024 20:38
 Operator : RC/JU
 Sample : P3314-01MSD
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ206MSD

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/24/2024
 Supervised By :mohammad ahmed 07/30/2024

Quant Time: Jul 24 01:16:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 09 16:58:40 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	90294	20.000	ng	-0.02
21) Naphthalene-d8	8.140	136	354478	20.000	ng	-0.01
39) Acenaphthene-d10	9.892	164	172809	20.000	ng	-0.01
64) Phenanthrene-d10	11.375	188	234304	20.000	ng	-0.01
76) Chrysene-d12	14.016	240	147778	20.000	ng	-0.01
86) Perylene-d12	15.474	264	121596	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	453173	79.486	ng	0.00
7) Phenol-d6	6.493	99	608263	80.200	ng	-0.01
23) Nitrobenzene-d5	7.422	82	389552	53.956	ng	-0.01
42) 2,4,6-Tribromophenol	10.681	330	115117	71.277	ng	-0.01
45) 2-Fluorobiphenyl	9.210	172	639151	57.806	ng	-0.02
79) Terphenyl-d14	12.951	244	438601	48.352	ng	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.710	88	98922	39.176	ng	97
3) Pyridine	3.457	79	257436	41.342	ng	96
4) n-Nitrosodimethylamine	3.422	42	188497	46.501	ng	94
6) Aniline	6.516	93	275350	38.761	ng	# 34
8) 2-Chlorophenol	6.645	128	294185	48.799	ng	99
9) Benzaldehyde	6.404	77	24104	5.938	ng	100
10) Phenol	6.510	94	387728	48.163	ng	88
11) bis(2-Chloroethyl)ether	6.593	93	287962	47.509	ng	98
12) 1,3-Dichlorobenzene	6.793	146	296692	43.784	ng	98
13) 1,4-Dichlorobenzene	6.875	146	304521	44.254	ng	98
14) 1,2-Dichlorobenzene	7.022	146	288769	45.057	ng	97
15) Benzyl Alcohol	6.998	79	280824	49.338	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.128	45	511687	42.929	ng	80
17) 2-Methylphenol	7.116	107	232501	47.058	ng	# 91
18) Hexachloroethane	7.363	117	114854	45.402	ng	95
19) n-Nitroso-di-n-propyla...	7.269	70	224543	47.914	ng	93
20) 3+4-Methylphenols	7.269	107	290700	46.724	ng	90
22) Acetophenone	7.269	105	374259	43.472	ng	99
24) Nitrobenzene	7.440	77	335417	45.716	ng	96
25) Isophorone	7.681	82	596663	48.123	ng	99
26) 2-Nitrophenol	7.757	139	156389	49.732	ng	97
27) 2,4-Dimethylphenol	7.792	122	197333	51.057	ng	100
28) bis(2-Chloroethoxy)met...	7.887	93	353695	47.732	ng	98
29) 2,4-Dichlorophenol	7.998	162	228905	45.608	ng	98
30) 1,2,4-Trichlorobenzene	8.075	180	251380	42.995	ng	98
31) Naphthalene	8.157	128	839958	45.561	ng	100
32) Benzoic acid	7.928	122	134768	36.560	ng	93
33) 4-Chloroaniline	8.210	127	149803	23.159	ng	98
34) Hexachlorobutadiene	8.269	225	147328	40.920	ng	99
35) Caprolactam	8.587	113	68478m	42.282	ng	
36) 4-Chloro-3-methylphenol	8.698	107	257386	45.808	ng	99
37) 2-Methylnaphthalene	8.851	142	520752	43.894	ng	100
38) 1-Methylnaphthalene	8.951	142	483195	41.732	ng	99
40) 1,2,4,5-Tetrachloroben...	9.016	216	216156	44.880	ng	99
41) Hexachlorocyclopentadiene	8.998	237	109471	70.026	ng	99
43) 2,4,6-Trichlorophenol	9.128	196	147046	47.863	ng	99

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44) 2,4,5-Trichlorophenol	9.181	196	153125	45.316	ng	95
46) 1,1'-Biphenyl	9.316	154	603790	46.104	ng	100
47) 2-Chloronaphthalene	9.339	162	471374	47.644	ng	99
48) 2-Nitroaniline	9.439	65	174497	50.897	ng	97
49) Acenaphthylene	9.757	152	719448	51.218	ng	99
50) Dimethylphthalate	9.616	163	566120	50.663	ng	99
51) 2,6-Dinitrotoluene	9.681	165	120595	46.873	ng	90
52) Acenaphthene	9.928	154	452347	48.446	ng	99
53) 3-Nitroaniline	9.845	138	92979	35.260	ng	91
54) 2,4-Dinitrophenol	9.957	184	86102	63.386	ng	95
55) Dibenzofuran	10.098	168	634449	46.847	ng	100
56) 4-Nitrophenol	10.022	139	140979	72.457	ng	98
57) 2,4-Dinitrotoluene	10.086	165	149487	44.906	ng	92
58) Fluorene	10.439	166	483369	45.107	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.216	232	112276	42.102	ng	96
60) Diethylphthalate	10.310	149	527672	48.984	ng	99
61) 4-Chlorophenyl-phenyle...	10.428	204	239377	44.486	ng	92
62) 4-Nitroaniline	10.463	138	99768	37.756	ng	99
63) Azobenzene	10.592	77	548038	47.319	ng	98
65) 4,6-Dinitro-2-methylph...	10.492	198	66607	49.233	ng	83
66) n-Nitrosodiphenylamine	10.551	169	411536	58.624	ng	98
67) 4-Bromophenyl-phenylether	10.922	248	130911	54.360	ng	98
68) Hexachlorobenzene	10.986	284	135325	50.582	ng	95
69) Atrazine	11.075	200	114143	49.043	ng	98
70) Pentachlorophenol	11.180	266	121367	76.666	ng	96
71) Phenanthrene	11.398	178	625219	52.818	ng	99
72) Anthracene	11.451	178	592080	50.388	ng	100
73) Carbazole	11.610	167	470625	44.568	ng	100
74) Di-n-butylphthalate	11.933	149	652907	53.641	ng	100
75) Fluoranthene	12.586	202	574886	47.595	ng	98
77) Benzidine	12.710	184	150569	40.384	ng	99
78) Pyrene	12.816	202	566132	37.738	ng	99
80) Butylbenzylphthalate	13.427	149	218869	48.500	ng	95
81) Benzo(a)anthracene	14.004	228	527464	53.194	ng	99
82) 3,3'-Dichlorobenzidine	13.963	252	131351	51.408	ng	# 97
83) Chrysene	14.039	228	477292	50.881	ng	99
84) Bis(2-ethylhexyl)phtha...	13.986	149	312676	64.966	ng	99
85) Di-n-octyl phthalate	14.598	149	584944	76.934	ng	# 99
87) Indeno(1,2,3-cd)pyrene	16.939	276	303469	37.155	ng	99
88) Benzo(b)fluoranthene	15.051	252	445356	65.035	ng	99
89) Benzo(k)fluoranthene	15.074	252	376989	56.446	ng	99
90) Benzo(a)pyrene	15.410	252	342453	56.927	ng	100
91) Dibenzo(a,h)anthracene	16.951	278	233361	34.928	ng	99
92) Benzo(g,h,i)perylene	17.380	276	233711	32.988	ng	98

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

