

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031122\
 Data File : BF127298.D
 Acq On : 11 Mar 2022 12:24
 Operator : CG\JU
 Sample : N1645-01
 Misc : LOD-MDL-SOIL 8ppm
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-SOIL-01-QT1-2022

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: Mar 11 13:01:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 12:42:13 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							03/14/2022
1) 1,4-Dichlorobenzene-d4	6.869	152	91358	20.000	ng	0.00	Supervised By :Jagrut padhyay
21) Naphthalene-d8	8.152	136	328738	20.000	ng	# 0.00	
39) Acenaphthene-d10	9.910	164	205652	20.000	ng	0.00	
64) Phenanthrene-d10	11.404	188	353077	20.000	ng	# 0.00	
76) Chrysene-d12	14.051	240	302699	20.000	ng	# 0.00	
86) Perylene-d12	15.533	264	229516	20.000	ng	0.00	03/14/2022
System Monitoring Compounds							
5) 2-Fluorophenol	5.493	112	704848	124.027	ng	0.00	
7) Phenol-d6	6.505	99	1000990	128.039	ng	0.00	
23) Nitrobenzene-d5	7.434	82	570402	71.657	ng	0.00	
42) 2,4,6-Tribromophenol	10.710	330	274074	123.801	ng	0.00	
45) 2-Fluorobiphenyl	9.228	172	1003270	68.972	ng	0.00	
79) Terphenyl-d14	12.992	244	1137609	64.715	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	2.646	88	19757	6.705	ng	# 78	
3) Pyridine	3.475	79	53175	6.781	ng	# 74	
4) n-Nitrosodimethylamine	3.370	42	31547	7.261	ng	# 71	
6) Aniline	6.534	93	77712	8.318	ng	96	
8) 2-Chlorophenol	6.658	128	42243	7.185	ng	95	
9) Benzaldehyde	6.422	77	43269	9.495	ng	86	
10) Phenol	6.516	94	63510	7.395	ng	78	
11) bis(2-Chloroethyl)ether	6.605	93	48509	7.606	ng	92	
12) 1,3-Dichlorobenzene	6.810	146	49802	7.546	ng	98	
13) 1,4-Dichlorobenzene	6.887	146	49571	7.487	ng	97	
14) 1,2-Dichlorobenzene	7.040	146	46375	7.443	ng	97	
15) Benzyl Alcohol	7.010	79	45374	7.743	ng	# 89	
16) 2,2'-oxybis(1-Chloropr...	7.140	45	86973	7.267	ng	99	
17) 2-Methylphenol	7.116	107	37257	7.454	ng	# 92	
18) Hexachloroethane	7.375	117	18323	7.205	ng	# 81	
19) n-Nitroso-di-n-propyla...	7.275	70	35658	7.309	ng	# 86	
20) 3+4-Methylphenols	7.275	107	44268	6.830	ng	# 43	
22) Acetophenone	7.275	105	67998	7.555	ng	# 88	
24) Nitrobenzene	7.452	77	57675	7.661	ng	91	
25) Isophorone	7.687	82	96608	7.548	ng	99	
26) 2-Nitrophenol	7.769	139	20747	6.654	ng	# 64	
27) 2,4-Dimethylphenol	7.799	122	39728	8.488	ng	94	
28) bis(2-Chloroethoxy)met...	7.899	93	61463	7.579	ng	98	
29) 2,4-Dichlorophenol	8.010	162	37226	6.945	ng	98	
30) 1,2,4-Trichlorobenzene	8.093	180	44066	7.211	ng	95	
31) Naphthalene	8.175	128	129544	7.387	ng	99	
32) Benzoic acid	7.893	122	12001m	4.132	ng		
33) 4-Chloroaniline	8.222	127	55836	8.214	ng	# 94	
34) Hexachlorobutadiene	8.281	225	29046	7.128	ng	97	
35) Caprolactam	8.563	113	11225	7.010	ng	88	
36) 4-Chloro-3-methylphenol	8.704	107	41793	7.194	ng	97	
37) 2-Methylnaphthalene	8.863	142	88021	7.682	ng	96	
38) 1-Methylnaphthalene	8.963	142	87364	7.931	ng	97	
40) 1,2,4,5-Tetrachloroben...	9.028	216	46690	7.172	ng	# 93	
41) Hexachlorocyclopentadiene	9.010	237	10607	4.636	ng	98	
43) 2,4,6-Trichlorophenol	9.146	196	26335	6.409	ng	96	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.187	196	25322	5.807	ng	91	03/14/2022
46) 1,1'-Biphenyl	9.328	154	113678	7.324	ng	98	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	9.357	162	85352	7.064	ng	95	
48) 2-Nitroaniline	9.451	65	32457	6.982	ng	85	
49) Acenaphthylene	9.769	152	134872	7.294	ng	99	
50) Dimethylphthalate	9.622	163	99136	6.727	ng	# 99	
51) 2,6-Dinitrotoluene	9.693	165	20323	6.818	ng	90	03/14/2022
52) Acenaphthene	9.940	154	80129	7.270	ng	97	
53) 3-Nitroaniline	9.863	138	22703	6.989	ng	# 95	
54) 2,4-Dinitrophenol	9.981	184	4168	2.838	ng	# 90	
55) Dibenzofuran	10.116	168	119233	6.963	ng	99	
56) 4-Nitrophenol	10.034	139	10663	5.401	ng	# 78	
57) 2,4-Dinitrotoluene	10.099	165	28248	7.202	ng	94	
58) Fluorene	10.457	166	94597	7.125	ng	100	
59) 2,3,4,6-Tetrachlorophenol	10.234	232	21921	6.049	ng	# 82	
60) Diethylphthalate	10.322	149	94180	7.006	ng	97	
61) 4-Chlorophenyl-phenylether	10.446	204	50749	7.053	ng	89	
62) 4-Nitroaniline	10.475	138	22088	6.824	ng	# 76	
63) Azobenzene	10.610	77	104447	6.760	ng	90	
65) 4,6-Dinitro-2-methylph...	10.510	198	11508	5.286	ng	89	
66) n-Nitrosodiphenylamine	10.569	169	80934	7.381	ng	98	
67) 4-Bromophenyl-phenylether	10.940	248	30189	7.093	ng	93	
68) Hexachlorobenzene	11.004	284	32201	6.942	ng	# 90	
69) Atrazine	11.093	200	30674	8.106	ng	95	
70) Pentachlorophenol	11.210	266	8116m	4.334	ng		
71) Phenanthrene	11.428	178	134778	7.231	ng	98	
72) Anthracene	11.475	178	141538	7.668	ng	98	
73) Carbazole	11.634	167	123046	7.079	ng	98	
74) Di-n-butylphthalate	11.957	149	142406	7.004	ng	100	
75) Fluoranthene	12.616	202	155836	7.478	ng	92	
77) Benzidine	12.740	184	66696	8.664	ng	98	
78) Pyrene	12.845	202	158789	6.735	ng	98	
80) Butylbenzylphthalate	13.457	149	58170	6.411	ng	# 95	
81) Benzo(a)anthracene	14.039	228	142707	7.002	ng	100	
82) 3,3'-Dichlorobenzidine	13.998	252	45671	7.207	ng	# 96	
83) Chrysene	14.075	228	130520	6.885	ng	98	
84) Bis(2-ethylhexyl)phtha...	14.010	149	82923	6.789	ng	# 100	
85) Di-n-octyl phthalate	14.628	149	134848	7.294	ng	96	
87) Indeno(1,2,3-cd)pyrene	17.039	276	83481	6.038	ng	# 85	
88) Benzo(b)fluoranthene	15.092	252	113573	7.539	ng	# 93	
89) Benzo(k)fluoranthene	15.122	252	106991	7.553	ng	# 95	
90) Benzo(a)pyrene	15.469	252	99364	8.376	ng	# 93	
91) Dibenzo(a,h)anthracene	17.051	278	73613	6.397	ng	# 91	
92) Benzo(g,h,i)perylene	17.492	276	69000	6.147	ng	# 84	

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

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