

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061019\
 Data File : BE099897.D
 Acq On : 10 Jun 2019 19:20
 Operator : JU/SJ
 Sample : K2241-03
 Misc : MDL-SOIL-0.12PPM
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-MDL-SOIL-03-QT2-2019

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:39:15 AM

Quant Time: Jun 11 07:16:40 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Jun 10 14:57:20 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	612	0.40	ng	0.00
7) Naphthalene-d8	10.61	136	2170	0.40	ng	0.01
13) Acenaphthene-d10	14.47	164	1723	0.40	ng	0.00
19) Phenanthrene-d10	17.23	188	5953	0.40	ng	0.01
27) Chrysene-d12	21.40	240	7947	0.40	ng	0.00
34) Perylene-d12	23.93	264	9218	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.36	112	1471	0.89	ng	0.00
5) Phenol-d6	6.96	99	1821	0.86	ng	0.00
8) Nitrobenzene-d5	8.98	82	1258	0.61	ng	0.01
11) 2-Methylnaphthalene-d10	12.22	152	1527	0.40	ng	0.01
14) 2,4,6-Tribromophenol	15.98	330	929	0.73	ng	0.01
15) 2-Fluorobiphenyl	13.10	172	4511	0.69	ng	0.00
25) Fluoranthene-d10	19.25	212	32092	0.37	ng	0.00
29) Terphenyl-d14	19.85	244	12748	0.89	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.27	88	138	0.152	ng	# 85
3) n-Nitrosodimethylamine	3.64	42	76	0.146	ng	# 100
6) bis(2-Chloroethyl)ether	7.23	93	230	0.132	ng	# 61
9) Naphthalene	10.67	128	2803	0.120	ng	# 91
10) Hexachlorobutadiene	10.93	225	225	0.126	ng	# 98
12) 2-Methylnaphthalene	12.29	142	450	0.118	ng	# 94
16) Acenaphthylene	14.20	152	975	0.114	ng	# 96
17) Acenaphthene	14.54	154	523	0.112	ng	# 99
18) Fluorene	15.53	166	839	0.120	ng	# 96
20) 4-Bromophenyl-phenylether	16.42	248	388	0.115	ng	# 90
21) Hexachlorobenzene	16.52	284	420	0.122	ng	# 96
22) Pentachlorophenol	16.99	266	22m	0.313	ng	
23) Phenanthrene	17.27	178	1549	0.107	ng	# 100
24) Anthracene	17.36	178	1223	0.092	ng	# 98
26) Fluoranthene	19.28	202	2783	0.113	ng	# 99
28) Pyrene	19.65	202	2787	0.133	ng	# 100
30) Benzo(a)anthracene	21.39	228	2620	0.110	ng	# 95
31) Chrysene	21.44	228	3143	0.128	ng	# 96
32) Bis(2-ethylhexyl)phthalate	21.29	149	3954	0.116	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.63	276	3541	0.122	ng	# 98
35) Benzo(b)fluoranthene	23.15	252	2785	0.108	ng	# 86
36) Benzo(k)fluoranthene	23.20	252	3275	0.121	ng	# 90
37) Benzo(a)pyrene	23.82	252	3066	0.120	ng	# 83
38) Dibenzo(a,h)anthracene	26.66	278	2725	0.106	ng	# 79
39) Benzo(g,h,i)perylene	27.44	276	3129	0.118	ng	# 87

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

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