

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE111218\
 Data File : BE098194.D
 Acq On : 12 Nov 2018 17:36
 Operator : SJ/JU
 Sample : J5807-13DL 500X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-BNA-SoilDL

Quant Time: Nov 13 09:34:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 12 13:32:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1086	0.40	ng	0.01
7) Naphthalene-d8	10.67	136	5328	0.40	ng	0.00
13) Acenaphthene-d10	14.53	164	3448	0.40	ng	0.00
19) Phenanthrene-d10	17.28	188	8258	0.40	ng	0.00
27) Chrysene-d12	21.46	240	9150	0.40	ng	0.00
34) Perylene-d12	24.01	264	8782	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.45	112	768	0.25	ng	0.00
5) Phenol-d6	7.04	99	1382	0.28	ng	0.00
8) Nitrobenzene-d5	9.03	82	1016	0.19	ng	0.01
11) 2-Methylnaphthalene-d10	12.22	152	8	0.00	ng	-0.04
14) 2,4,6-Tribromophenol	16.03	330	366	0.24	ng	0.01
15) 2-Fluorobiphenyl	13.15	172	2705	0.19	ng	0.00
25) Fluoranthene-d10	19.31	212	24	0.00	ng	0.00
29) Terphenyl-d14	19.91	244	4067	0.20	ng	0.00

Target Compounds				Qvalue		
6) bis(2-Chloroethyl)ether	7.29	93	398	0.100	ng	96
20) 4-Bromophenyl-phenylether	16.46	248	2672	0.531	ng	95
21) Hexachlorobenzene	16.58	284	3531	0.675	ng	98
22) Pentachlorophenol	16.94	266	540	0.543	ng	98
23) Phenanthrene	17.32	178	10009	0.487	ng	99
28) Pyrene	19.70	202	15271	0.606	ng	99
30) Benzo(a)anthracene	21.45	228	5469	0.210	ng	99
32) Bis(2-ethylhexyl)phthalate	21.37	149	39869	0.738	ng	100
33) Indeno(1,2,3-cd)pyrene	26.70	276	12736	0.526	ng	# 93
36) Benzo(k)fluoranthene	23.28	252	13012	0.495	ng	100
38) Dibenzo(a,h)anthracene	26.73	278	5278	0.259	ng	94

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

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