

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122018\
 Data File : BE098542.D
 Acq On : 20 Dec 2018 17:00
 Operator : SJ/JU
 Sample : PB115586BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB115586BS

Quant Time: Dec 21 05:03:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	840	0.40	ng	-0.01
7) Naphthalene-d8	10.64	136	3784	0.40	ng	-0.03
13) Acenaphthene-d10	14.51	164	2684	0.40	ng	-0.01
19) Phenanthrene-d10	17.26	188	7390	0.40	ng	-0.03
27) Chrysene-d12	21.45	240	7960	0.40	ng	-0.01
34) Perylene-d12	23.99	264	7127	0.40	ng	-0.02

System Monitoring Compounds						
4) 2-Fluorophenol	5.44	112	842	0.43	ng	0.00
5) Phenol-d6	7.03	99	1166	0.42	ng	-0.01
8) Nitrobenzene-d5	9.02	82	1276	0.37	ng	0.00
11) 2-Methylnaphthalene-d10	12.25	152	2658	0.43	ng	-0.01
14) 2,4,6-Tribromophenol	16.01	330	486	0.49	ng	-0.01
15) 2-Fluorobiphenyl	13.13	172	4392	0.39	ng	-0.01
25) Fluoranthene-d10	19.29	212	40221	0.42	ng	-0.02
29) Terphenyl-d14	19.89	244	8008	0.41	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.36	88	619	0.426	ng	98
3) n-Nitrosodimethylamine	3.69	42	390	0.298	ng	# 100
6) bis(2-Chloroethyl)ether	7.27	93	913	0.345	ng	95
9) Naphthalene	10.69	128	16237	0.426	ng	99
10) Hexachlorobutadiene	10.98	225	1084	0.447	ng	98
12) 2-Methylnaphthalene	12.33	142	2618	0.423	ng	97
16) Acenaphthylene	14.23	152	4738	0.377	ng	99
17) Acenaphthene	14.57	154	3128	0.414	ng	98
18) Fluorene	15.55	166	4293	0.425	ng	99
20) 4-Bromophenyl-phenylether	16.45	248	1592	0.400	ng	92
21) Hexachlorobenzene	16.57	284	1891	0.442	ng	96
22) Pentachlorophenol	16.93	266	468	0.629	ng	94
23) Phenanthrene	17.30	178	7442	0.414	ng	100
24) Anthracene	17.40	178	5939	0.371	ng	99
26) Fluoranthene	19.32	202	9927	0.418	ng	99
28) Pyrene	19.68	202	10357	0.415	ng	99
30) Benzo(a)anthracene	21.43	228	9154	0.404	ng	98
31) Chrysene	21.49	228	9199	0.387	ng	98
32) Bis(2-ethylhexyl)phthalate	21.35	149	15188	0.330	ng	99
33) Indeno(1,2,3-cd)pyrene	26.66	276	9303	0.394	ng	98
35) Benzo(b)fluoranthene	23.21	252	7850	0.403	ng	99
36) Benzo(k)fluoranthene	23.25	252	9104	0.401	ng	98
37) Benzo(a)pyrene	23.87	252	7840	0.390	ng	97
38) Dibenzo(a,h)anthracene	26.68	278	7690	0.406	ng	99
39) Benzo(g,h,i)perylene	27.48	276	8021	0.420	ng	98

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

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