

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121118\
 Data File : BE098356.D
 Acq On : 12 Dec 2018 6:02
 Operator : SJ/JU
 Sample : J6314-22
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Dec 12 17:26:56 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	7.86	152	1059	0.40	ng	0.00
7) Naphthalene-d8	10.67	136	4530	0.40	ng	0.00
13) Acenaphthene-d10	14.53	164	2943	0.40	ng	0.00
19) Phenanthrene-d10	17.28	188	7915	0.40	ng	0.00
27) Chrysene-d12	21.46	240	9472	0.40	ng	0.00
34) Perylene-d12	24.01	264	9088	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.45	112	474	0.19	ng	0.00
5) Phenol-d6	7.04	99	387	0.11	ng	0.00
8) Nitrobenzene-d5	9.03	82	1974	0.48	ng	0.01
11) 2-Methylnaphthalene-d10	12.27	152	3301	0.45	ng	0.00
14) 2,4,6-Tribromophenol	16.02	330	577	0.53	ng	0.00
15) 2-Fluorobiphenyl	13.15	172	6686	0.54	ng	0.00
25) Fluoranthene-d10	19.31	212	49574	0.49	ng	0.00
29) Terphenyl-d14	19.90	244	13985	0.61	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.37	88	209	0.114	ng	# 45
3) n-Nitrosodimethylamine	3.63	42	160	0.097	ng	# 100
22) Pentachlorophenol	16.97	266	23	0.049	ng	# 76
26) Fluoranthene	19.34	202	543	0.021	ng	# 88
30) Benzo(a)anthracene	21.45	228	594	0.022	ng	# 72
31) Chrysene	21.50	228	589	0.021	ng	# 72
32) Bis(2-ethylhexyl)phthalate	21.37	149	5258	0.096	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.71	276	624	0.022	ng	# 78
35) Benzo(b)fluoranthene	23.23	252	589	0.024	ng	# 4
37) Benzo(a)pyrene	23.90	252	544	0.021	ng	# 5
38) Dibenzo(a,h)anthracene	26.73	278	490	0.020	ng	# 21
39) Benzo(g,h,i)perylene	27.53	276	562	0.023	ng	# 50

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

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