

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE111618\
 Data File : BE098237.D
 Acq On : 16 Nov 2018 19:45
 Operator : SJ/JU
 Sample : J5957-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 W-1

Manual Integrations
 APPROVED

Sohil
 11/19/2018 3:44:03 PM

Quant Time: Nov 16 22:59:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 16 19:05:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	1264	0.40	ng	0.00
7) Naphthalene-d8	10.67	136	6155	0.40	ng	0.00
13) Acenaphthene-d10	14.53	164	4120	0.40	ng	0.00
19) Phenanthrene-d10	17.28	188	9614	0.40	ng	0.00
27) Chrysene-d12	21.46	240	10458	0.40	ng	0.00
34) Perylene-d12	24.01	264	9900	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.46	112	397	0.12	ng	0.00
5) Phenol-d6	7.04	99	399	0.07	ng	0.00
8) Nitrobenzene-d5	9.03	82	1693	0.29	ng	0.01
11) 2-Methylnaphthalene-d10	12.27	152	3021	0.28	ng	0.00
14) 2,4,6-Tribromophenol	16.02	330	448	0.34	ng	-0.01
15) 2-Fluorobiphenyl	13.15	172	5020	0.28	ng	0.00
25) Fluoranthene-d10	19.31	212	35814	0.31	ng	0.00
29) Terphenyl-d14	19.91	244	5958	0.25	ng	0.00

Target Compounds					Qvalue	
16) Acenaphthylene	14.26	152	3268	0.167	ng	99
17) Acenaphthene	14.60	154	1122	0.096	ng	96
18) Fluorene	15.58	166	1162	0.078	ng	# 97
23) Phenanthrene	17.31	178	4669	0.189	ng	97
24) Anthracene	17.41	178	3328	0.148	ng	99
26) Fluoranthene	19.34	202	6458	0.218	ng	# 95
28) Pyrene	19.70	202	10671	0.349	ng	# 96
30) Benzo(a)anthracene	21.45	228	2515	0.083	ng	# 65
31) Chrysene	21.50	228	1871	0.060	ng	# 76
32) Bis(2-ethylhexyl)phthalate	21.37	149	18896	0.401	ng	99
33) Indeno(1,2,3-cd)pyrene	26.70	276	7100	0.266	ng	# 93
35) Benzo(b)fluoranthene	23.23	252	4975	0.178	ng	# 81
36) Benzo(k)fluoranthene	23.28	252	1503m	0.048	ng	
37) Benzo(a)pyrene	23.90	252	3777	0.138	ng	# 80
38) Dibenzo(a,h)anthracene	26.72	278	1394	0.063	ng	# 65
39) Benzo(g,h,i)perylene	27.53	276	14065	0.593	ng	97

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

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