

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE110118\
 Data File : BE098149.D
 Acq On : 1 Nov 2018 16:47
 Operator : SJ/JU
 Sample : J5665-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 50

Instrument :
 BNA_E
 ClientSampleId :
 TWP-04

Quant Time: Nov 02 13:17:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:00:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	60148	20.00	ng	-0.01
7) Naphthalene-d8	10.67	136	282714	20.00	ng	0.00
13) Acenaphthene-d10	14.53	164	192211	20.00	ng	-0.01
19) Phenanthrene-d10	17.28	188	451556	20.00	ng	0.00
27) Chrysene-d12	21.48	240	506620	20.00	ng	0.01
34) Perylene-d12	24.02	264	518087	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.46	112	154717	43.33	ng	0.00
5) Phenol-d6	7.04	99	166473	28.31	ng	-0.01
8) Nitrobenzene-d5	9.03	82	360657	59.48	ng	0.00
11) 2-Methylnaphthalene-d10	12.27	152	50	0.01	ng	0.00
14) 2,4,6-Tribromophenol	16.03	330	174063	95.62	ng	0.00
15) 2-Fluorobiphenyl	13.16	172	777603	49.20	ng	0.00
25) Fluoranthene-d10	19.33	212	43	0.00	ng	0.02
29) Terphenyl-d14	19.91	244	1094219	46.44	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Naphthalene	10.72	128	2661	0.047	ng	# 93
12) 2-Methylnaphthalene	12.34	142	622	0.062	ng	# 87
23) Phenanthrene	17.33	178	2312	0.103	ng	# 98
24) Anthracene	17.42	178	476	0.022	ng	# 93
26) Fluoranthene	19.34	202	3005	0.107	ng	# 98
28) Pyrene	19.71	202	2802	0.101	ng	# 96
30) Benzo(a)anthracene	21.45	228	4259	0.144	ng	# 92
33) Indeno(1,2,3-cd)pyrene	26.71	276	1067	0.035	ng	# 89
35) Benzo(b)fluoranthene	23.24	252	3727	0.128	ng	# 88
37) Benzo(a)pyrene	23.90	252	1019	0.036	ng	# 60
39) Benzo(a,h,i)perylene	27.53	276	1049	0.040	ng	# 74

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

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