

Data Path : Z:\HPCHEM1\BNA E\DATA\BE121417\
 Data File : BE094986.D
 Acq On : 14 Dec 2017 20:48
 Operator : SJ/JU
 Sample : I6797-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 YT-MW-07

Manual Integrations
 APPROVED

Sohil
 12/15/2017 7:54:42 PM

Quant Time: Dec 15 08:23:49 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 14 18:42:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	6663	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	14691	0.40	ng	0.00
13) Acenaphthene-d10	15.06	164	9608	0.40	ng	0.00
19) Phenanthrene-d10	17.77	188	19863	0.40	ng	0.00
27) Chrysene-d12	21.87	240	24229	0.40	ng	0.00
34) Perylene-d12	24.55	264	18079	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.95	112	1634	0.16	ng	0.00
5) Phenol-d6	7.59	99	1357	0.09	ng	0.00
8) Nitrobenzene-d5	9.65	82	4693	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	12.86	152	8022	0.33	ng	0.00
14) 2,4,6-Tribromophenol	16.53	330	1236	0.55	ng	0.00
15) 2-Fluorobiphenyl	13.70	172	15429	0.36	ng	0.00
25) Fluoranthene-d10	19.75	212	90778	0.35	ng	0.00
29) Terphenyl-d14	20.31	244	20732	0.45	ng	0.00
Target Compounds						
9) Naphthalene	11.38	128	180992	1.049	ng	99
12) 2-Methylnaphthalene	12.93	142	2182	0.051	ng	# 80
17) Acenaphthene	15.12	154	16585	0.487	ng	# 85
18) Fluorene	16.08	166	16328	0.386	ng	# 82
22) Pentachlorophenol	17.42	266	211	0.151	ng	89
23) Phenanthrene	17.82	178	6906	0.112	ng	# 61
30) Benzo(a)anthracene	21.85	228	2626m	0.037	ng	
31) Chrysene	21.92	228	2392m	0.030	ng	
33) Indeno(1,2,3-cd)pyrene	27.40	276	1985	0.031	ng	# 83
35) Benzo(b)fluoranthene	23.72	252	2499	0.038	ng	# 70
38) Dibenzo(a,h)anthracene	27.43	278	1606	0.029	ng	# 69
39) Benzo(g,h,i)perylene	28.29	276	1688	0.031	ng	# 72

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

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