

Data Path : Z:\HPCHEM1\BNA E\DATA\BE121417\
 Data File : BE094990.D
 Acq On : 14 Dec 2017 23:08
 Operator : SJ/JU
 Sample : I6876-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 KY038PS034-171211

Manual Integrations
 APPROVED

Sohil
 12/15/2017 7:54:52 PM

Quant Time: Dec 15 08:28:35 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	6727	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	14459	0.40	ng	0.00
13) Acenaphthene-d10	15.06	164	7978	0.40	ng	0.00
19) Phenanthrene-d10	17.77	188	18963	0.40	ng	0.00
27) Chrysene-d12	21.87	240	21540	0.40	ng	0.00
34) Perylene-d12	24.55	264	17983	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.95	112	1636	0.16	ng	0.00
5) Phenol-d6	7.59	99	1403	0.09	ng	0.00
8) Nitrobenzene-d5	9.65	82	5112	0.43	ng	0.00
11) 2-Methylnaphthalene-d10	12.86	152	8837	0.37	ng	0.00
14) 2,4,6-Tribromophenol	16.53	330	745	0.40	ng	0.00
15) 2-Fluorobiphenyl	13.70	172	14360	0.41	ng	0.00
25) Fluoranthene-d10	19.75	212	95838	0.39	ng	0.00
29) Terphenyl-d14	20.31	244	19402	0.48	ng	0.00
Target Compounds						
30) Benzo(a)anthracene	21.85	228	2444	0.039	ng	Qvalue # 65
31) Chrysene	21.90	228	2279m	0.032	ng	
33) Indeno(1,2,3-cd)pyrene	27.41	276	2352	0.041	ng	100
35) Benzo(b)fluoranthene	23.71	252	2417	0.037	ng	# 82
36) Benzo(k)fluoranthene	23.77	252	2301	0.034	ng	# 60
37) Benzo(a)pyrene	24.42	252	2036	0.035	ng	# 52
38) Dibenzo(a,h)anthracene	27.43	278	1868	0.034	ng	# 79
39) Benzo(g,h,i)perylene	28.29	276	2034	0.038	ng	# 82

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

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