

Data Path : Z:\HPCHEM1\BNA E\DATA\BE121417\
 Data File : BE094987.D
 Acq On : 14 Dec 2017 21:24
 Operator : SJ/JU
 Sample : I6797-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EB-2017-12-07

Manual Integrations
 APPROVED

Sohil
 12/15/2017 7:54:45 PM

Quant Time: Dec 15 08:25:11 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	5615	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	12487	0.40	ng	0.00
13) Acenaphthene-d10	15.06	164	6893	0.40	ng	0.00
19) Phenanthrene-d10	17.77	188	16525	0.40	ng	0.00
27) Chrysene-d12	21.87	240	20154	0.40	ng	0.00
34) Perylene-d12	24.54	264	16434	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.96	112	1116	0.13	ng	0.00
5) Phenol-d6	7.59	99	953	0.07	ng	0.00
8) Nitrobenzene-d5	9.65	82	4342	0.42	ng	0.00
11) 2-Methylnaphthalene-d10	12.86	152	6776	0.33	ng	0.00
14) 2,4,6-Tribromophenol	16.53	330	373	0.23	ng	0.00
15) 2-Fluorobiphenyl	13.70	172	12851	0.42	ng	0.00
25) Fluoranthene-d10	19.75	212	76128	0.36	ng	0.00
29) Terphenyl-d14	20.31	244	17864	0.47	ng	0.00
Target Compounds						
30) Benzo(a)anthracene	21.85	228	2186	0.037	ng	Qvalue # 67
31) Chrysene	21.90	228	2093m	0.032	ng	
33) Indeno(1,2,3-cd)pyrene	27.40	276	2214	0.041	ng	100
35) Benzo(b)fluoranthene	23.72	252	2287	0.038	ng	# 87
36) Benzo(k)fluoranthene	23.77	252	2059	0.034	ng	# 66
37) Benzo(a)pyrene	24.42	252	1867	0.035	ng	# 56
38) Dibenzo(a,h)anthracene	27.43	278	1687	0.034	ng	# 79
39) Benzo(g,h,i)perylene	28.29	276	1920	0.039	ng	# 83

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

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