

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE112817\
 Data File : BE094880.D
 Acq On : 28 Nov 2017 12:51
 Operator : SJ/JU
 Sample : I6514-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB2

Manual Integrations
 APPROVED

Sohil
 11/29/2017 5:14:57 PM

Quant Time: Nov 29 01:24:17 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 29 01:10:42 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	577	0.40	ng/ul	-0.05
2) Naphthalene-d8	10.70	136	3778	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	2593	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	6092	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	6005	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	6058	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1898	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	5509	0.26	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	246	0.027	ng/ul#	50
5) 2-Methylnaphthalene	12.38	142	251	0.040	ng/ul	96
12) Phenanthrene	17.27	178	2372	0.145	ng/ul#	87
15) Fluoranthene	19.30	202	4546	0.212	ng/ul	84
17) Pyrene	19.66	202	4190	0.223	ng/ul#	82
18) Benzo(a)anthracene	21.41	228	2393	0.131	ng/ul#	92
19) Chrysene	21.47	228	2767	0.152	ng/ul	92
21) Benzo(b)fluoranthene	23.18	252	3337m	0.185	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	1311m	0.068	ng/ul	
23) Benzo(a)pyrene	23.84	252	2153	0.120	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.69	276	1512	0.077	ng/ul#	78
26) Benzo(g,h,i)perylene	27.54	276	935	0.057	ng/ul#	67

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

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