

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093578.D
 Acq On : 25 Jul 2017 15:35
 Operator : SJ/JU
 Sample : I4209-15
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B18

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:55:45 PM

Quant Time: Jul 25 18:09:42 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 18:00:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2490	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12892	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	8465	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	19915m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	21241	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	17752m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5155	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	17458	0.28	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	2092	0.067	ng/ul#	96
12) Phenanthrene	17.29	178	1400	0.026	ng/ul#	91
15) Fluoranthene	19.30	202	2448	0.036	ng/ul	87
17) Pyrene	19.67	202	2940	0.047	ng/ul#	90
18) Benzo(a)anthracene	21.40	228	1932	0.031	ng/ul#	89
19) Chrysene	21.45	228	1978	0.033	ng/ul#	89
21) Benzo(b)fluoranthene	23.14	252	3182	0.058	ng/ul#	74
23) Benzo(a)pyrene	23.80	252	2789m	0.053	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	1832m	0.032	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	1767m	0.036	ng/ul	

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

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