

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072517\
 Data File : BE093590.D
 Acq On : 25 Jul 2017 22:57
 Operator : SJ/JU
 Sample : I4209-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0B15

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:56:09 PM

Quant Time: Jul 26 10:39:16 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 06:10:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2750	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	14426	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	9663	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	22338m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23887	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	20604m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7841	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	23229	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	1786	0.051	ng/ul#	90
7) Acenaphthylene	14.25	152	1027	0.025	ng/ul#	89
12) Phenanthrene	17.29	178	1373	0.023	ng/ul#	91
15) Fluoranthene	19.30	202	6805	0.089	ng/ul	85
17) Pyrene	19.67	202	6809m	0.098	ng/ul	
18) Benzo(a)anthracene	21.40	228	4435	0.064	ng/ul#	87
19) Chrysene	21.45	228	3203	0.048	ng/ul	94
21) Benzo(b)fluoranthene	23.14	252	6100	0.096	ng/ul	91
22) Benzo(k)fluoranthene	23.19	252	2391m	0.038	ng/ul	
23) Benzo(a)pyrene	23.80	252	5068m	0.084	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	2853m	0.043	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	2554m	0.045	ng/ul	

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

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