

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013118\
 Data File : BE095337.D
 Acq On : 1 Feb 2018 3:07
 Operator : SJ/JU
 Sample : J1245-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B78

Manual Integrations
 APPROVED

Sohil
 2/1/2018 5:59:00 PM

Quant Time: Feb 01 06:08:26 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 04:05:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2310	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7286	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4887	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9792m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9801	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9018m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4795	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13109	0.35	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	17.36	266	2374	1.252	ng/ul	95
12) Phenanthrene	17.75	178	1846	0.056	ng/ul#	89
13) Anthracene	17.85	178	611	0.020	ng/ul#	75
15) Fluoranthene	19.72	202	6429	0.151	ng/ul	84
17) Pyrene	20.08	202	5177m	0.157	ng/ul	
18) Benzo(a)anthracene	21.79	228	2365	0.077	ng/ul#	90
19) Chrysene	21.84	228	2623	0.082	ng/ul#	89
21) Benzo(b)fluoranthene	23.63	252	2991m	0.094	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	1187m	0.041	ng/ul	
23) Benzo(a)pyrene	24.32	252	1438m	0.051	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.25	276	1166m	0.037	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	1077m	0.041	ng/ul	

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

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