

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095512.D
 Acq On : 6 Feb 2018 3:56
 Operator : SJ/JU
 Sample : J1315-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C82

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:31:39 PM

Quant Time: Feb 06 06:20:50 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 04:22:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1821	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5225	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3779	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8265	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7161	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6610m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2636	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8078	0.26	ng/ul	0.00
Target Compounds					Qvalue	
7) Acenaphthylene	14.75	152	1964	0.108	ng/ul#	75
11) Pentachlorophenol	17.36	266	1623	1.014	ng/ul	98
12) Phenanthrene	17.75	178	6277m	0.226	ng/ul	
13) Anthracene	17.82	178	5418m	0.213	ng/ul	
15) Fluoranthene	19.72	202	6482	0.180	ng/ul	78
17) Pyrene	20.08	202	14834m	0.617	ng/ul	
18) Benzo(a)anthracene	21.78	228	3957	0.177	ng/ul#	2
19) Chrysene	21.81	228	6086	0.262	ng/ul#	45
21) Benzo(b)fluoranthene	23.62	252	9649m	0.414	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	2302m	0.108	ng/ul	
23) Benzo(a)pyrene	24.32	252	6301m	0.303	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	3615m	0.158	ng/ul	
25) Dibenzo(a,h)anthracene	27.24	278	799m	0.043	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	3073m	0.159	ng/ul	

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

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