

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020918\
 Data File : BE095639.D
 Acq On : 10 Feb 2018 10:06
 Operator : SJ/JU
 Sample : J1355-01
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6D26

Manual Integrations
 APPROVED

Sohil
 2/12/2018 11:40:28 AM

Quant Time: Feb 12 01:36:20 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 10 02:59:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1954	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5854	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3899	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8964	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	8162	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7522m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3061	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8405	0.25	ng/ul	0.00
Target Compounds				Qvalue		
5) 2-Methylnaphthalene	12.88	142	293	0.027	ng/ul	99
11) Pentachlorophenol	17.36	266	2509	1.445	ng/ul	99
12) Phenanthrene	17.75	178	932m	0.031	ng/ul	
15) Fluoranthene	19.72	202	3586	0.092	ng/ul	84
17) Pyrene	20.07	202	3501m	0.128	ng/ul	
18) Benzo(a)anthracene	21.78	228	1368	0.054	ng/ul#	89
19) Chrysene	21.83	228	1265	0.048	ng/ul#	86
21) Benzo(b)fluoranthene	23.61	252	1971m	0.074	ng/ul	
23) Benzo(a)pyrene	24.31	252	626m	0.026	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.22	276	524m	0.020	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	492m	0.022	ng/ul	

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

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