

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020918\
 Data File : BE095628.D
 Acq On : 10 Feb 2018 3:35
 Operator : SJ/JU
 Sample : J1339-03
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C93

Manual Integrations
 APPROVED

Sohil
 2/12/2018 11:39:37 AM

Quant Time: Feb 12 00:30:34 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	1834	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5408	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3393	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8257	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7564	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7902	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	1111	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	3113	0.10	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	600	0.037	ng/ul#	74
11) Pentachlorophenol	17.36	266	9233	5.773	ng/ul	98
12) Phenanthrene	17.75	178	4350	0.157	ng/ul	93
13) Anthracene	17.84	178	1939	0.076	ng/ul	98
15) Fluoranthene	19.72	202	24486	0.681	ng/ul	83
17) Pyrene	20.07	202	24108m	0.950	ng/ul	
18) Benzo(a)anthracene	21.78	228	12116	0.512	ng/ul#	85
19) Chrysene	21.83	228	9456	0.385	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	14447m	0.518	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	3280m	0.129	ng/ul	
23) Benzo(a)pyrene	24.31	252	5085m	0.204	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.22	276	3567	0.131	ng/ul#	73
25) Dibenzo(a,h)anthracene	27.25	278	861	0.039	ng/ul#	60
26) Benzo(g,h,i)perylene	28.11	276	3041	0.132	ng/ul	96

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

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