

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020918\
 Data File : BE095626.D
 Acq On : 10 Feb 2018 2:23
 Operator : SJ/JU
 Sample : J1339-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6D40

Manual Integrations
 APPROVED

Sohil
 2/12/2018 11:39:31 AM

Quant Time: Feb 10 03:17:37 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	8.42	152	1920	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5560	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3606	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8240	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7679	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7185m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2252	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	6191	0.20	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	665	0.038	ng/ul#	79
8) Acenaphthene	15.07	153	807	0.060	ng/ul#	92
9) Fluorene	16.02	166	2477	0.178	ng/ul	94
11) Pentachlorophenol	17.36	266	65484	41.030	ng/ul	98
12) Phenanthrene	17.75	178	61974m	2.236	ng/ul	
13) Anthracene	17.84	178	6923	0.273	ng/ul	95
15) Fluoranthene	19.72	202	64304	1.792	ng/ul	82
17) Pyrene	20.07	202	47241m	1.833	ng/ul	
18) Benzo(a)anthracene	21.78	228	13879	0.578	ng/ul#	85
19) Chrysene	21.84	228	10093	0.405	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	9989m	0.394	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	2459m	0.106	ng/ul	
23) Benzo(a)pyrene	24.31	252	3901m	0.172	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.23	276	2173m	0.088	ng/ul	
25) Dibenzo(a,h)anthracene	27.24	278	546m	0.027	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	2044m	0.097	ng/ul	

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

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