

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095467.D
 Acq On : 4 Feb 2018 8:52
 Operator : SJ/JU
 Sample : J1315-19
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C82

Manual Integrations
 APPROVED

Sohil
 2/5/2018 7:01:01 PM

Quant Time: Feb 05 06:55:26 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 04:46:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1805	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5392	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	4100	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7916	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7155	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6850m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3124	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8532	0.28	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	2791	0.141	ng/ul#	82
11) Pentachlorophenol	17.36	266	1779	1.160	ng/ul	98
12) Phenanthrene	17.75	178	7771m	0.292	ng/ul	
13) Anthracene	17.82	178	5728m	0.235	ng/ul	
15) Fluoranthene	19.72	202	6859	0.199	ng/ul	77
17) Pyrene	20.08	202	16680	0.695	ng/ul#	84
18) Benzo(a)anthracene	21.79	228	4776	0.213	ng/ul#	38
19) Chrysene	21.82	228	5615	0.242	ng/ul#	38
21) Benzo(b)fluoranthene	23.62	252	10115m	0.419	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	2614m	0.118	ng/ul	
23) Benzo(a)pyrene	24.32	252	6445m	0.299	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	3800m	0.161	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	3210m	0.160	ng/ul	

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

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