

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095473.D
 Acq On : 4 Feb 2018 12:27
 Operator : SJ/JU
 Sample : J1315-03
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C60

Manual Integrations
 APPROVED

Sohil
 2/5/2018 7:01:21 PM

Quant Time: Feb 05 07:42:24 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	8.43	152	1896	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5795	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3850	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8531	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7986	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7366m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3000	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	7727	0.24	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	523	0.028	ng/ul#	73
11) Pentachlorophenol	17.36	266	4088	2.474	ng/ul	98
12) Phenanthrene	17.75	178	1562m	0.054	ng/ul	
13) Anthracene	17.85	178	2137m	0.081	ng/ul	
15) Fluoranthene	19.72	202	14107	0.380	ng/ul	85
17) Pyrene	20.08	202	13910	0.519	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	3591	0.144	ng/ul#	90
19) Chrysene	21.84	228	4288	0.165	ng/ul	94
21) Benzo(b)fluoranthene	23.62	252	6581m	0.253	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	1763m	0.074	ng/ul	
23) Benzo(a)pyrene	24.32	252	1639m	0.071	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.23	276	1734m	0.068	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	1447m	0.067	ng/ul	

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

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