

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095242.D
 Acq On : 29 Jan 2018 12:21
 Operator : SJ/JU
 Sample : J1233-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B03

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:31:53 PM

Quant Time: Jan 30 04:53:37 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 30 04:43:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1957	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5846	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4018	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7951m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	8434	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7393m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3981	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	11565	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.32	128	434	0.027	ng/ul#	55
5) 2-Methylnaphthalene	12.89	142	1561	0.143	ng/ul	100
7) Acenaphthylene	14.74	152	869	0.045	ng/ul#	66
8) Acenaphthene	15.08	153	7666	0.511	ng/ul	98
9) Fluorene	16.02	166	9227	0.597	ng/ul#	70
11) Pentachlorophenol	17.36	266	9598	6.232	ng/ul	95
12) Phenanthrene	17.75	178	40684	1.521	ng/ul#	86
13) Anthracene	17.84	178	11545	0.472	ng/ul#	92
15) Fluoranthene	19.73	202	47022	1.358	ng/ul	81
17) Pyrene	20.08	202	30924	1.093	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	7116	0.270	ng/ul#	87
19) Chrysene	21.84	228	5076	0.185	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	7173m	0.275	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	1705m	0.071	ng/ul	
23) Benzo(a)pyrene	24.32	252	2072m	0.089	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1787m	0.070	ng/ul	
25) Dibenzo(a,h)anthracene	27.27	278	435m	0.021	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	2090m	0.097	ng/ul	

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

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