

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095252.D
 Acq On : 29 Jan 2018 18:20
 Operator : SJ/JU
 Sample : J1233-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B15

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:32:10 PM

Quant Time: Jan 30 06:02:51 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	2576	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	8064	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5739	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10717m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10678	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9450m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5609	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	14972	0.37	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	11.31	128	79178	3.506	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	35964	2.385	ng/ul	100
7) Acenaphthylene	14.74	152	1603	0.058	ng/ul#	53
8) Acenaphthene	15.08	153	42768	1.997	ng/ul	94
9) Fluorene	16.02	166	40807	1.847	ng/ul#	73
11) Pentachlorophenol	17.36	266	169869	81.835	ng/ul	95
12) Phenanthrene	17.75	178	145657	4.041	ng/ul#	86
13) Anthracene	17.85	178	27056	0.820	ng/ul#	92
15) Fluoranthene	19.72	202	64443	1.381	ng/ul	84
17) Pyrene	20.08	202	46978	1.311	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	11471	0.344	ng/ul#	88
19) Chrysene	21.84	228	7284	0.210	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	6015m	0.180	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	1782m	0.058	ng/ul	
23) Benzo(a)pyrene	24.32	252	3064m	0.103	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	796m	0.024	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	1068m	0.039	ng/ul	

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

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