

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097481.D
 Acq On : 13 Sep 2018 9:53
 Operator : SJ/JU
 Sample : J4792-08DL 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z39DL

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:43:08 PM

Quant Time: Sep 13 11:37:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 13 07:48:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1023	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5837	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3658	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	8658	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9956	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	10006	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	327	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	1005	0.04	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.17	128	7007	0.492	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	2382	0.265	ng/ul	98
7) Acenaphthylene	13.73	152	4195	0.315	ng/ul	99
8) Acenaphthene	14.07	153	2190	0.191	ng/ul	96
9) Fluorene	15.07	166	5651	0.430	ng/ul	93
12) Phenanthrene	16.80	178	43721	1.962	ng/ul#	89
13) Anthracene	16.90	178	10176	0.543	ng/ul#	93
15) Fluoranthene	18.83	202	39190	1.369	ng/ul	84
17) Pyrene	19.20	202	36183	1.351	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	19652	0.768	ng/ul#	86
19) Chrysene	21.01	228	17657m	0.599	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	28459m	1.062	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	8504m	0.285	ng/ul	
23) Benzo(a)pyrene	23.11	252	21470	0.866	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.51	276	19393	0.604	ng/ul#	72
25) Dibenzo(a,h)anthracene	25.52	278	4639	0.170	ng/ul#	65
26) Benzo(g,h,i)perylene	26.21	276	17325	0.632	ng/ul	93

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

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