

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097450.D
 Acq On : 12 Sep 2018 12:11
 Operator : SJ/JU
 Sample : J4792-06MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YZ4MS

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:14 PM

Quant Time: Sep 13 01:30:29 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 01:16:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	921	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5259	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3562	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	9163	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10775	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11491	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1599	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	6450	0.21	ng/ul	0.00
Target Compounds				Qvalue		
8) Acenaphthene	14.07	153	2024	0.181	ng/ul	95
11) Pentachlorophenol	16.43	266	108	0.120	ng/ul	99
12) Phenanthrene	16.80	178	1661	0.070	ng/ul#	89
13) Anthracene	16.90	178	521	0.026	ng/ul#	57
15) Fluoranthene	18.83	202	3817	0.126	ng/ul	82
17) Pyrene	19.20	202	28391	0.979	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	48782	1.762	ng/ul#	79
19) Chrysene	21.00	228	65783	2.063	ng/ul#	74
21) Benzo(b)fluoranthene	22.54	252	11092m	0.360	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1901m	0.055	ng/ul	
23) Benzo(a)pyrene	23.10	252	11793	0.414	ng/ul#	23
24) Indeno(1,2,3-cd)pyrene	25.50	276	3470	0.094	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.51	278	7208	0.231	ng/ul#	22
26) Benzo(g,h,i)perylene	26.20	276	2907	0.092	ng/ul#	29

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

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