

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097451.D
 Acq On : 12 Sep 2018 12:49
 Operator : SJ/JU
 Sample : J4792-07MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YZ4MSD

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:15 PM

Quant Time: Sep 13 01:31:56 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	7.38	152	910	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5285	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3640	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	9367	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10497	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11205	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1780	0.22	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	6877	0.22	ng/ul	0.00
Target Compounds				Qvalue		
8) Acenaphthene	14.07	153	2118	0.186	ng/ul	96
11) Pentachlorophenol	16.43	266	159	0.172	ng/ul	94
12) Phenanthrene	16.80	178	1682	0.070	ng/ul#	89
13) Anthracene	16.89	178	557	0.027	ng/ul#	48
15) Fluoranthene	18.84	202	4145	0.134	ng/ul	80
17) Pyrene	19.20	202	30506	1.080	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	52897	1.962	ng/ul#	78
19) Chrysene	21.01	228	70836	2.280	ng/ul#	75
21) Benzo(b)fluoranthene	22.54	252	11795m	0.393	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	2182m	0.065	ng/ul	
23) Benzo(a)pyrene	23.11	252	13220	0.476	ng/ul#	22
24) Indeno(1,2,3-cd)pyrene	25.50	276	3723	0.103	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.51	278	8016	0.263	ng/ul#	21
26) Benzo(g,h,i)perylene	26.20	276	2877	0.094	ng/ul#	30

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

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