

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097417.D
 Acq On : 7 Sep 2018 9:59
 Operator : SJ/JU
 Sample : J4688-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YW1

Quant Time: Sep 07 23:50:48 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 : Fri Sep 07 23:49:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	787	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	3860	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2349	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	7126	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9338	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9211	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1987	0.33	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	9305	0.40	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.17	128	2720	0.289	ng/ul#	95
5) 2-Methylnaphthalene	11.80	142	1639	0.276	ng/ul	98
7) Acenaphthylene	13.72	152	2689	0.314	ng/ul	98
9) Fluorene	15.07	166	3221	0.382	ng/ul	94
11) Pentachlorophenol	16.44	266	138	0.197	ng/ul	92
12) Phenanthrene	16.80	178	12644	0.689	ng/ul#	89
13) Anthracene	16.89	178	1941	0.126	ng/ul	93
15) Fluoranthene	18.83	202	11161	0.474	ng/ul	84
17) Pyrene	19.20	202	13126	0.522	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	16112	0.672	ng/ul#	86
19) Chrysene	21.01	228	11168	0.404	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	7074	0.287	ng/ul	85
22) Benzo(k)fluoranthene	22.57	252	13670	0.497	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.48	276	11156	0.377	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.51	278	10260	0.409	ng/ul#	91
26) Benzo(g,h,i)perylene	26.18	276	14836	0.588	ng/ul#	90

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

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