

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098114.D
 Acq On : 31 Oct 2018 12:39
 Operator : SJ/JU
 Sample : J5635-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A40C5

Quant Time: Nov 01 01:44:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 01:42:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	933	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4842	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3345	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	7907	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	10450	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	8942	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1930	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	6779	0.26	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.73	128	2582	0.223	ng/ul	97
5) 2-Methylnaphthalene	12.34	142	1933	0.221	ng/ul	99
7) Acenaphthylene	14.25	152	7222	0.547	ng/ul	98
9) Fluorene	15.59	166	3321	0.280	ng/ul	99
12) Phenanthrene	17.33	178	10634	0.583	ng/ul	99
13) Anthracene	17.42	178	3816	0.212	ng/ul	98
15) Fluoranthene	19.35	202	9246	0.397	ng/ul	95
17) Pyrene	19.71	202	12010	0.492	ng/ul	99
18) Benzo(a)anthracene	21.45	228	19817	0.645	ng/ul	98
19) Chrysene	21.51	228	8628	0.363	ng/ul	99
21) Benzo(b)fluoranthene	23.23	252	6890	0.278	ng/ul	94
22) Benzo(k)fluoranthene	23.29	252	10731	0.442	ng/ul#	98
24) Indeno(1,2,3-cd)pyrene	26.72	276	9304	0.341	ng/ul#	90
25) Dibenzo(a,h)anthracene	26.74	278	8658	0.380	ng/ul#	95
26) Benzo(g,h,i)perylene	27.54	276	12214	0.531	ng/ul	98

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

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