

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM070816\
 Data File : BM006312.D
 Acq On : 07 Jul 2016 10:26
 Operator : UM/SJ
 Sample : H3811-27
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A4TX7

Quant Time: Jul 07 13:46:43 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM070716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 07 05:51:09 2016
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.650	152	4169	0.40	ng/ul	0.00
4) Naphthalene-d8	10.427	136	16392	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.297	164	9449	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.039	188	20695	0.40	ng/ul	0.00
18) Chrysene-d12	21.242	240	13533	0.40	ng/ul	0.00
22) Perylene-d12	23.451	264	11711	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.135	96	83	0.05	ng/uL	-0.03
6) 2-Methylnaphthalene-d10	12.018	152	6391	0.33	ng/ul	0.00
16) Fluoranthene-d10	19.072	212	19759	0.45	ng/ul	-0.01
Target Compounds						
					Ovalue	
5) Naphthalene	10.476	128	17931	0.47	ng/ul	99
7) 2-Methylnaphthalene	12.091	142	13213	0.52	ng/ul	99
9) Acenaphthylene	14.006	152	16550	0.52	ng/ul	93
10) Acenaphthene	14.348	153	19908	0.63	ng/ul	90
13) Pentachlorophenol	16.696	266	3085	0.97	ng/ul	96
14) Phenanthrene	17.176	178	26007	0.45	ng/ul	99
15) Anthracene	17.176	178	26342	0.50	ng/ul	97
17) Fluoranthene	19.108	202	1546	0.02	ng/ul#	1
19) Pyrene	19.471	202	49652	0.88	ng/ul	95
20) Benzo(a)anthracene	21.273	228	30269	0.69	ng/ul#	92
21) Chrysene	21.273	228	30269	0.70	ng/ul	95
23) Benzo(b)fluoranthene	22.784	252	33923	0.81	ng/ul	98
24) Benzo(k)fluoranthene	22.826	252	47594	1.22	ng/ul	96
26) Indeno(1,2,3-cd)pyrene	25.672	276	7210	0.17	ng/ul#	51
27) Dibenzo(a,h)anthracene	25.682	278	28748	0.85	ng/ul	97

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

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