

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061319\
 Data File : BN006306.D
 Acq On : 13 Jun 2019 11:56
 Operator : JU/SJ
 Sample : K3213-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8H1

Quant Time: Jun 13 12:44:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 13 11:20:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	3514	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	13701	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	7174	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	13661	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	10997	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	12498	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6909	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	14466	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	9403	0.249	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	6536	0.248	ng/ul	99
7) Acenaphthylene	14.01	152	28945	0.726	ng/ul	99
8) Acenaphthene	14.36	153	8803	0.318	ng/ul	98
9) Fluorene	15.36	166	9956	0.308	ng/ul	97
11) Pentachlorophenol	16.72	266	412	0.119	ng/ul	90
12) Phenanthrene	17.09	178	29735	0.678	ng/ul	100
13) Anthracene	17.19	178	9427	0.225	ng/ul	97
15) Fluoranthene	19.12	202	21488	0.444	ng/ul	98
17) Pyrene	19.49	202	25085	0.433	ng/ul	98
18) Benzo(a)anthracene	21.25	228	35606	0.732	ng/ul	98
19) Chrysene	21.30	228	22430	0.491	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	18168	0.349	ng/ul	97
23) Benzo(a)pyrene	23.40	252	24102	0.497	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.74	276	22638	0.431	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.75	278	19108	0.455	ng/ul	97
26) Benzo(a,h,i)perylene	26.43	276	23091	0.525	ng/ul	97

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

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