

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\
 Data File : BN006216.D
 Acq On : 10 Jun 2019 11:15
 Operator : JU/SJ
 Sample : K3016-27
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51H5

Quant Time: Jun 10 11:50:44 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 : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5086	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	19830	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9972	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	22015	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	17980	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	20297	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.02	152	8937	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	21786	0.37	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.48	128	54667	0.999	ng/ul	99
7) Acenaphthylene	14.01	152	28966	0.523	ng/ul	100
8) Acenaphthene	14.36	153	22911	0.596	ng/ul	100
9) Fluorene	15.35	166	20093	0.447	ng/ul	97
11) Pentachlorophenol	16.71	266	3032	0.542	ng/ul	97
12) Phenanthrene	17.09	178	35978	0.509	ng/ul	99
13) Anthracene	17.18	178	34183	0.506	ng/ul	100
15) Fluoranthene	19.12	202	43051	0.552	ng/ul	98
17) Pyrene	19.49	202	59718	0.631	ng/ul	98
18) Benzo(a)anthracene	21.25	228	42372	0.533	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	42390	0.501	ng/ul#	62
23) Benzo(a)pyrene	23.40	252	35470	0.450	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.74	276	63532	0.745	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.75	278	48047	0.704	ng/ul	93
26) Benzo(g,h,i)perylene	26.41	276	56204	0.787	ng/ul	95

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

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