

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062419\
 Data File : BN006524.D
 Acq On : 24 Jun 2019 10:30
 Operator : HP/JU
 Sample : K3360-24
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41T2

Quant Time: Jun 25 11:07:31 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 : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5858	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	22015	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.28	164	12061	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.04	188	25479	0.40	ng/ul	-0.02
16) Chrysene-d12	21.27	240	16528	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	16074	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.01	152	8985	0.27	ng/ul	-0.02
14) Fluoranthene-d10	19.08	212	19747	0.29	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	13987	0.230	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	18783	0.444	ng/ul	100
7) Acenaphthylene	14.00	152	14031	0.209	ng/ul	99
8) Acenaphthene	14.34	153	14321	0.308	ng/ul	100
9) Fluorene	15.34	166	21845	0.401	ng/ul	98
11) Pentachlorophenol	16.71	266	2500	0.386	ng/ul	99
12) Phenanthrene	17.08	178	26162	0.320	ng/ul	99
13) Anthracene	17.17	178	30792	0.394	ng/ul	99
15) Fluoranthene	19.11	202	28673	0.317	ng/ul	97
17) Pyrene	19.48	202	28782	0.331	ng/ul	99
18) Benzo(a)anthracene	21.25	228	28130	0.385	ng/ul	98
19) Chrysene	21.30	228	28565	0.416	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	33259	0.496	ng/ul	89
22) Benzo(k)fluoranthene	22.88	252	18051	0.276	ng/ul	98
23) Benzo(a)pyrene	23.41	252	23796	0.381	ng/ul	94
25) Dibenzo(a,h)anthracene	25.75	278	32679	0.605	ng/ul	93
26) Benzo(g,h,i)perylene	26.43	276	14285	0.252	ng/ul	96

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

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