

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072219\
 Data File : BN006943.D
 Acq On : 22 Jul 2019 12:25
 Operator : HP/JU
 Sample : K3770-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2H33

Quant Time: Jul 22 13:39:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 19 23:29:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1900	0.40	ng/ul	0.01
2) Naphthalene-d8	10.38	136	7183	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.25	164	3818	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	7579	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	5898	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	6151	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3551	0.32	ng/ul	0.02
14) Fluoranthene-d10	19.06	212	7863	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	17830	0.865	ng/ul	99
7) Acenaphthylene	13.97	152	13657	0.693	ng/ul	99
8) Acenaphthene	14.31	153	11200	0.737	ng/ul	99
9) Fluorene	15.32	166	13757	0.826	ng/ul	100
11) Pentachlorophenol	16.71	266	1347	1.193	ng/ul	97
12) Phenanthrene	17.06	178	17850	0.732	ng/ul	99
13) Anthracene	17.17	178	7719	0.366	ng/ul	98
17) Pyrene	19.45	202	46248	1.716	ng/ul	96
18) Benzo(a)anthracene	21.23	228	14947	0.617	ng/ul	100
19) Chrysene	21.28	228	29927	1.107	ng/ul	99
22) Benzo(k)fluoranthene	22.86	252	12481	0.473	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.71	276	19539	0.782	ng/ul#	85
26) Benzo(g,h,i)perylene	26.39	276	23158	1.108	ng/ul	97

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

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