

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN032919\
 Data File : BN005031.D
 Acq On : 29 Mar 2019 13:30
 Operator : JU/SJ
 Sample : K1825-05DL 25X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-BN-WPDL

Quant Time: Mar 30 00:18:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN032819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 29 03:42:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3069	0.40	ng	0.00
7) Naphthalene-d8	10.44	136	11772	0.40	ng	0.00
13) Acenaphthene-d10	14.30	164	7145	0.40	ng	0.00
19) Phenanthrene-d10	17.07	188	16852	0.40	ng	0.01
27) Chrysene-d12	21.26	240	21509	0.40	ng	0.00
34) Perylene-d12	23.49	264	19477	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	42367	5.53	ng	0.00
5) Phenol-d6	6.83	99	53798	5.93	ng	-0.02
8) Nitrobenzene-d5	8.82	82	29671	4.35	ng	-0.01
11) 2-Methylnaphthalene-d10	12.04	152	2	0.00	ng	0.00
14) 2,4,6-Tribromophenol	15.80	330	23899	5.96	ng	-0.01
15) 2-Fluorobiphenyl	12.92	172	112566	4.17	ng	0.00
25) Fluoranthene-d10	19.09	212	16	0.00	ng	0.00
29) Terphenyl-d14	19.69	244	188160	4.26	ng	0.00

Target Compounds				Qvalue		
3) n-Nitrosodimethylamine	3.51	42	5844	2.902	ng	# 92
9) Naphthalene	10.49	128	103566	3.500	ng	98
16) Acenaphthylene	14.02	152	118314	3.508	ng	99
17) Acenaphthene	14.37	154	13398	0.626	ng	99
20) 4-Bromophenyl-phenylether	16.25	248	24271	2.515	ng	92
26) Fluoranthene	19.12	202	373931	6.472	ng	100
30) Benzo(a)anthracene	21.24	228	373002	6.150	ng	99
31) Chrysene	21.29	228	346473	5.299	ng	98
32) Bis(2-ethylhexyl)phthalate	21.15	149	185108	7.288	ng	99
37) Benzo(a)pyrene	23.39	252	131013	2.360	ng	95
39) Benzo(g,h,i)perylene	26.42	276	288245	5.152	ng	98

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

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