

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
 Data File : BP001338.D
 Acq On : 20 Dec 2019 12:01
 Operator : JU
 Sample : K6325-04DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RT-3425DL

Quant Time: Dec 20 13:09:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 03:20:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	67277	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	250028	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	144107	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	273858	20.00	ng	0.00
76) Chrysene-d12	19.25	240	187106	20.00	ng	0.00
87) Perylene-d12	21.02	264	182808	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.20	112	32572	7.82	ng	0.00
7) Phenol-d6	7.74	99	55495	10.22	ng	-0.01
23) Nitrobenzene-d5	9.18	82	39805	6.11	ng	-0.01
42) 2,4,6-Tribromophenol	14.49	330	10762	5.97	ng	0.00
45) 2-Fluorobiphenyl	12.11	172	65378	5.74	ng	0.00
79) Terphenyl-d14	17.88	244	71947	6.59	ng	0.00

Target Compounds

						Qvalue
31) Naphthalene	10.37	128	94229	6.904	ng	98
37) 2-Methylnaphthalene	11.50	142	60585	6.430	ng	99
38) 1-Methylnaphthalene	11.66	142	27604	3.116	ng	97
52) Acenaphthene	13.25	154	95046	11.069	ng	99
55) Dibenzofuran	13.53	168	92630	6.885	ng	99
58) Fluorene	14.10	166	90105	8.374	ng	97
71) Phenanthrene	15.64	178	416672	26.668	ng	100
72) Anthracene	15.71	178	54151	3.501	ng	97
73) Carbazole	15.96	167	32789	2.378	ng	99
75) Fluoranthene	17.35	202	237916	13.618	ng	99
78) Pyrene	17.65	202	157869	10.867	ng	97
81) Benzo(a)anthracene	19.23	228	31116	2.282	ng	95
83) Chrysene	19.27	228	31797	2.416	ng	94

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

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