

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070720\
 Data File : BP002663.D
 Acq On : 07 Jul 2020 18:05
 Operator : CG/JU
 Sample : L3208-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RT-4743

Quant Time: Jul 08 01:49:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	237899	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	915337	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	538886	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	547198	20.00	ng	0.00
76) Chrysene-d12	21.09	240	938772	20.00	ng	0.00
86) Perylene-d12	23.28	264	1269456	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	1193876	84.97	ng	0.00
7) Phenol-d6	6.76	99	1817695	92.68	ng	0.00
23) Nitrobenzene-d5	8.71	82	1152928	64.90	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	419065	63.94	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	2741583	74.88	ng	0.00
79) Terphenyl-d14	19.57	244	1876517	43.15	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.68	163	231910	5.479	ng	# 99
71) Phenanthrene	17.03	178	133179	4.524	ng	# 44
75) Fluoranthene	19.02	202	180447	5.063	ng	# 85
78) Pyrene	19.37	202	185422	2.902	ng	# 90
81) Benzo(a)anthracene	21.07	228	168833	2.751	ng	# 97
83) Chrysene	21.13	228	192945	3.278	ng	# 95
88) Benzo(b)fluoranthene	22.63	252	317679	3.922	ng	# 92
90) Benzo(a)pyrene	23.19	252	196606	2.645	ng	# 91

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

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