

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
 Data File : BP003160.D
 Acq On : 28 Aug 2020 20:32
 Operator : CG/JU
 Sample : L3837-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 789UMR-TP07-1218-01

Quant Time: Aug 29 02:08:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 15:32:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	250970	20.00	ng	0.00
21) Naphthalene-d8	10.44	136	1012086	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	620539	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1352398	20.00	ng	-0.01
76) Chrysene-d12	21.15	240	1420523	20.00	ng	-0.01
86) Perylene-d12	23.38	264	1324242	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	1117486	78.72	ng	0.00
7) Phenol-d6	6.85	99	1335032	75.03	ng	0.00
23) Nitrobenzene-d5	8.80	82	792812	56.47	ng	0.00
42) 2,4,6-Tribromophenol	15.80	330	669351	79.75	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	2364676	55.81	ng	0.00
79) Terphenyl-d14	19.63	244	3437971	53.26	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.76	163	197268	4.169	ng	99
71) Phenanthrene	17.09	178	192725	2.630	ng	99
75) Fluoranthene	19.07	202	508957	6.009	ng	99
78) Pyrene	19.42	202	409444	4.441	ng	100
81) Benzo(a)anthracene	21.13	228	209396	2.353	ng	100
83) Chrysene	21.18	228	245840	2.886	ng	98
88) Benzo(b)fluoranthene	22.71	252	325378	3.953	ng	99
90) Benzo(a)pyrene	23.28	252	189960	2.488	ng	98

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

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