

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082720\
 Data File : BP003138.D
 Acq On : 27 Aug 2020 18:40
 Operator : CG/JU
 Sample : L3816-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ALMASI-1

Quant Time: Aug 28 01:01:27 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	215728	20.00	ng	0.00
21) Naphthalene-d8	10.44	136	862364	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	533030	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1167472	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	1257688	20.00	ng	-0.01
86) Perylene-d12	23.38	264	1378529	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	857072	70.24	ng	0.00
7) Phenol-d6	6.84	99	1033486	67.57	ng	0.00
23) Nitrobenzene-d5	8.80	82	593060	49.58	ng	0.00
42) 2,4,6-Tribromophenol	15.80	330	514186	71.32	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	1842905	50.63	ng	0.00
79) Terphenyl-d14	19.63	244	2729768	47.76	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.76	163	173639	4.272	ng	99
71) Phenanthrene	17.10	178	290080	4.585	ng	99
75) Fluoranthene	19.07	202	312892	4.279	ng	100
78) Pyrene	19.42	202	280568	3.437	ng	99

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

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