

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092220\
 Data File : BP003342.D
 Acq On : 22 Sep 2020 16:55
 Operator : CG/JU
 Sample : L3869-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OR-02-091820

Quant Time: Sep 22 20:46:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 12:48:08 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	164489	20.00	ng	0.00
21) Naphthalene-d8	10.346	136	649532	20.00	ng	# 0.00
39) Acenaphthene-d10	14.216	164	402928	20.00	ng	0.00
64) Phenanthrene-d10	16.975	188	870144	20.00	ng	# 0.00
76) Chrysene-d12	21.069	240	942934	20.00	ng	0.00
86) Perylene-d12	23.263	264	909511	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	970499	103.02	ng	0.00
7) Phenol-d6	6.769	99	1149458	91.54	ng	0.00
23) Nitrobenzene-d5	8.722	82	889003	80.41	ng	0.00
42) 2,4,6-Tribromophenol	15.722	330	608387	99.00	ng	0.00
45) 2-Fluorobiphenyl	12.840	172	2043169	74.16	ng	0.00
79) Terphenyl-d14	19.557	244	3170258	70.05	ng	0.00
Target Compounds						
10) Phenol	6.793	94	37410	3.12	ng	79
50) Dimethylphthalate	13.681	163	153658	4.78	ng	99

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

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