

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090922\
 Data File : BP011666.D
 Acq On : 09 Sep 2022 16:47
 Operator : CG/JU
 Sample : N4539-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MBD-WC-S-MP-1-0-46

Quant Time: Sep 10 00:53:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP090222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 10 00:48:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	438355	20.000	ng	0.00
21) Naphthalene-d8	10.775	136	1703047	20.000	ng	# 0.00
39) Acenaphthene-d10	14.604	164	967688	20.000	ng	0.00
64) Phenanthrene-d10	17.357	188	1997085	20.000	ng	0.00
76) Chrysene-d12	21.433	240	1904597	20.000	ng	0.00
86) Perylene-d12	23.910	264	2041557	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	3465853	139.834	ng	0.00
7) Phenol-d6	7.116	99	4040207	123.191	ng	0.00
23) Nitrobenzene-d5	9.128	82	2520030	90.831	ng	0.00
42) 2,4,6-Tribromophenol	16.092	330	1828109	167.205	ng	0.00
45) 2-Fluorobiphenyl	13.228	172	6228248	92.038	ng	0.00
79) Terphenyl-d14	19.910	244	9505317	104.530	ng	0.00

Target Compounds	Qvalue

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

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