

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092421\
 Data File : BP007159.D
 Acq On : 24 Sep 2021 15:10
 Operator : CG/JU
 Sample : M3911-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SW-RR-01-(2.0)

Quant Time: Sep 24 17:06:27 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 23 11:17:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	242125	20.000	ng	0.00
21) Naphthalene-d8	10.687	136	927962	20.000	ng	# 0.00
39) Acenaphthene-d10	14.516	164	578639	20.000	ng	-0.01
64) Phenanthrene-d10	17.269	188	1219601	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1264590	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	1364581	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	923250	63.829	ng	0.00
7) Phenol-d6	7.057	99	681251	34.460	ng	0.00
23) Nitrobenzene-d5	9.051	82	1393589	87.458	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	1059569	141.243	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	3440567	85.183	ng	0.00
79) Terphenyl-d14	19.827	244	5319571	80.609	ng	0.00

Target Compounds	Qvalue

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

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