

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092221\
 Data File : BP007103.D
 Acq On : 22 Sep 2021 18:02
 Operator : CG/JU
 Sample : M3835-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SW-RR-02-(1.0)

Quant Time: Sep 22 18:29:22 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 16:50:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	273905	20.000	ng	0.00
21) Naphthalene-d8	10.698	136	1087582	20.000	ng	# 0.00
39) Acenaphthene-d10	14.528	164	686679	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	1406707	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1357475	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	1421861	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	834305	50.988	ng	0.00
7) Phenol-d6	7.063	99	627781	28.071	ng	0.00
23) Nitrobenzene-d5	9.057	82	1500140	80.327	ng	-0.01
42) 2,4,6-Tribromophenol	16.016	330	1135329	127.530	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	3837415	80.060	ng	0.00
79) Terphenyl-d14	19.833	244	5491848	77.525	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.087	94	57616	2.672	ng	93
20) 3+4-Methylphenols	8.663	107	365653	18.193	ng	87
31) Naphthalene	10.746	128	746831	13.455	ng	99

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

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