

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110421\
 Data File : BP007859.D
 Acq On : 04 Nov 2021 21:25
 Operator : CG/JU
 Sample : M4475-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB02A

Quant Time: Nov 07 11:19:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 19:18:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	186123	20.000	ng	-0.01
21) Naphthalene-d8	10.710	136	703531	20.000	ng	-0.02
39) Acenaphthene-d10	14.551	164	444682	20.000	ng	-0.01
64) Phenanthrene-d10	17.304	188	952170	20.000	ng	-0.02
76) Chrysene-d12	21.398	240	1054761	20.000	ng	-0.02
86) Perylene-d12	23.827	264	1188720	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	1302595	118.691	ng	0.00
7) Phenol-d6	7.087	99	1587754	99.955	ng	0.00
23) Nitrobenzene-d5	9.069	82	939454	71.731	ng	-0.02
42) 2,4,6-Tribromophenol	16.045	330	675809	106.489	ng	-0.01
45) 2-Fluorobiphenyl	13.175	172	2186169	71.594	ng	-0.02
79) Terphenyl-d14	19.869	244	3381401	65.280	ng	-0.01
Target Compounds						
32) Benzoic acid	9.981	122	45277	5.415	ng	Qvalue 95

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

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