

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112921\
 Data File : BP008139.D
 Acq On : 29 Nov 2021 13:42
 Operator : CG/JU
 Sample : M4843-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 L-1

Quant Time: Nov 29 14:08:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 25 01:13:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	77742	20.000	ng	-0.01
21) Naphthalene-d8	10.599	136	287725	20.000	ng	-0.01
39) Acenaphthene-d10	14.451	164	186293	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	419976	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	489825	20.000	ng	-0.01
86) Perylene-d12	23.698	264	530486	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	520793	107.860	ng	-0.01
7) Phenol-d6	6.981	99	673741	103.366	ng	-0.02
23) Nitrobenzene-d5	8.963	82	456453	72.140	ng	-0.01
42) 2,4,6-Tribromophenol	15.945	330	273229	114.747	ng	-0.01
45) 2-Fluorobiphenyl	13.069	172	905811	70.627	ng	-0.01
79) Terphenyl-d14	19.781	244	1528551	65.218	ng	0.00
Target Compounds						
10) Phenol	7.011	94	18136	2.708	ng	87

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

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