

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111222\
 Data File : BP012643.D
 Acq On : 13 Nov 2022 04:48
 Operator : CG/JU
 Sample : N5431-06
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DISSOLVED-20221176-COMP-11-MODIFIED-ELUTRIATE

Quant Time: Nov 14 03:27:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 13 22:44:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	352836	20.000	ng	0.00
21) Naphthalene-d8	10.504	136	1572244	20.000	ng	-0.01
39) Acenaphthene-d10	14.363	164	1122919	20.000	ng	-0.02
64) Phenanthrene-d10	17.128	188	2513083	20.000	ng	-0.02
76) Chrysene-d12	21.233	240	2525424	20.000	ng	-0.01
86) Perylene-d12	23.574	264	2305957	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	616399	32.530	ng	0.00
7) Phenol-d6	6.893	99	690137	26.239	ng	-0.02
23) Nitrobenzene-d5	8.881	82	1017168	35.892	ng	-0.01
42) 2,4,6-Tribromophenol	15.869	330	1000592	65.273	ng	-0.02
45) 2-Fluorobiphenyl	12.987	172	2473710	33.442	ng	-0.01
79) Terphenyl-d14	19.704	244	4316345	34.626	ng	-0.01

Target Compounds	Qvalue

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

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