

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011922\
 Data File : BP008859.D
 Acq On : 19 Jan 2022 18:09
 Operator : CG/JU
 Sample : N1154-10
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20220029-KEANSBURG-3-COMP-SPLP-LEACHATE

Quant Time: Jan 20 05:55:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 14 18:05:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	286983	20.000	ng	-0.01
21) Naphthalene-d8	10.646	136	1038405	20.000	ng	-0.01
39) Acenaphthene-d10	14.486	164	636308	20.000	ng	-0.01
64) Phenanthrene-d10	17.239	188	1205693	20.000	ng	-0.01
76) Chrysene-d12	21.333	240	1093827	20.000	ng	-0.01
86) Perylene-d12	23.745	264	1221659	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	506509	27.293	ng	0.00
7) Phenol-d6	7.022	99	347290	15.454	ng	-0.01
23) Nitrobenzene-d5	9.004	82	1332250	72.839	ng	-0.01
42) 2,4,6-Tribromophenol	15.981	330	606365	65.467	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	3451251	71.527	ng	-0.01
79) Terphenyl-d14	19.804	244	4363087	67.329	ng	-0.01
Target Compounds						
37) 2-Methylnaphthalene	12.310	142	110336	2.676	ng	Qvalue 97

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

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