

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034086.D
 Acq On : 20 Sep 2024 14:29
 Operator : JU/RC
 Sample : P3957-16
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE104D3-20240911

Quant Time: Sep 20 15:05:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.430	152	5239	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	14777	0.400	ng	0.00
13) Acenaphthene-d10	14.079	164	8682	0.400	ng	0.00
19) Phenanthrene-d10	16.828	188	18625	0.400	ng	#-0.01
29) Chrysene-d12	21.060	240	13213	0.400	ng	# 0.00
35) Perylene-d12	23.178	264	9916	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1941	0.131	ng	0.00
5) Phenol-d6	6.629	99	1171	0.068	ng	0.00
8) Nitrobenzene-d5	8.568	82	3550	0.314	ng	0.00
11) 2-Methylnaphthalene-d10	11.788	152	7823	0.358	ng	0.00
14) 2,4,6-Tribromophenol	15.586	330	1159	0.295	ng	0.00
15) 2-Fluorobiphenyl	12.690	172	12079	0.342	ng	-0.01
27) Fluoranthene-d10	18.878	212	20952	0.446	ng	0.00
31) Terphenyl-d14	19.501	244	13801	0.523	ng	0.00
Target Compounds						
25) Phenanthrene	16.877	178	1296	0.024	ng	# 96

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

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