

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102524\
 Data File : BN034729.D
 Acq On : 25 Oct 2024 18:53
 Operator : RC/JU
 Sample : P4550-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 VPB190-HYD-20241023

Quant Time: Oct 25 22:43:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 25 22:11:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.654	152	8808	0.400	ng	-0.06
7) Naphthalene-d8	10.426	136	25081	0.400	ng	#-0.05
13) Acenaphthene-d10	14.275	164	12081	0.400	ng	-0.05
19) Phenanthrene-d10	17.019	188	23223	0.400	ng	#-0.05
29) Chrysene-d12	21.214	240	12463	0.400	ng	-0.04
35) Perylene-d12	23.414	264	11231	0.400	ng	-0.05
System Monitoring Compounds						
4) 2-Fluorophenol	5.264	112	607	0.023	ng	-0.06
5) Phenol-d6	6.809	99	9	0.000	ng	-0.07
8) Nitrobenzene-d5	8.781	82	5660	0.269	ng	-0.06
11) 2-Methylnaphthalene-d10	12.014	152	679	0.020	ng	-0.06
14) 2,4,6-Tribromophenol	15.766	330	853	0.228	ng	-0.05
15) 2-Fluorobiphenyl	12.907	172	12439	0.259	ng	-0.04
27) Fluoranthene-d10	19.057	212	9426	0.180	ng	-0.04
31) Terphenyl-d14	19.665	244	10091	0.396	ng	-0.04
Target Compounds						Qvalue
2) 1,4-Dioxane	3.241	88	243	0.021	ng	# 73
3) n-Nitrosodimethylamine	3.523	42	538	0.032	ng	# 4
6) bis(2-Chloroethyl)ether	7.069	93	120285	4.288	ng	# 35
34) Bis(2-ethylhexyl)phtha...	21.160	149	1169	0.043	ng	# 98

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

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