

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034523.D
 Acq On : 10 Oct 2024 16:19
 Operator : RC/JU
 Sample : P4263-05
 Misc :
 ALS Vial : 99 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE114D1-20240930

Quant Time: Oct 10 17:06:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 26 14:44:02 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	4547	0.400	ng	-0.02
7) Naphthalene-d8	10.127	136	12709	0.400	ng	-0.02
13) Acenaphthene-d10	14.019	164	7204	0.400	ng	-0.03
19) Phenanthrene-d10	16.783	188	13646	0.400	ng	#-0.02
29) Chrysene-d12	21.008	240	8251	0.400	ng	0.00
35) Perylene-d12	23.116	264	11639	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1401	0.109	ng	-0.01
5) Phenol-d6	6.585	99	614	0.041	ng	-0.01
8) Nitrobenzene-d5	8.525	82	2097	0.216	ng	-0.01
11) 2-Methylnaphthalene-d10	11.731	152	5418	0.289	ng	-0.02
14) 2,4,6-Tribromophenol	15.542	330	816	0.250	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	7532	0.257	ng	-0.02
27) Fluoranthene-d10	18.829	212	12789	0.371	ng	-0.02
31) Terphenyl-d14	19.447	244	9291	0.564	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	18470	3.249	ng	99
6) bis(2-Chloroethyl)ether	6.816	93	549	0.041	ng	91

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

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