

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034095.D
 Acq On : 20 Sep 2024 19:55
 Operator : JU/RC
 Sample : P4003-07
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE105D2-20240912

Quant Time: Sep 21 01:20:19 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	4603	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	12880	0.400	ng	0.00
13) Acenaphthene-d10	14.072	164	7819	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	17235	0.400	ng	# 0.00
29) Chrysene-d12	21.053	240	13336	0.400	ng	0.00
35) Perylene-d12	23.180	264	11229	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1391	0.106	ng	0.00
5) Phenol-d6	6.629	99	771	0.051	ng	0.00
8) Nitrobenzene-d5	8.578	82	2684	0.272	ng	0.00
11) 2-Methylnaphthalene-d10	11.787	152	6091	0.320	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	921	0.260	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	9848	0.310	ng	0.00
27) Fluoranthene-d10	18.876	212	19357	0.445	ng	-0.01
31) Terphenyl-d14	19.498	244	13114	0.492	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.068	88	22250	3.866	ng	# 87
6) bis(2-Chloroethyl)ether	6.874	93	701	0.052	ng	96

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

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