

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034107.D
 Acq On : 21 Sep 2024 03:52
 Operator : JU/RC
 Sample : P4003-16
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE108D2-20240913

Quant Time: Sep 21 04:23:18 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	7.431	152	4883	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	13576	0.400	ng	0.00
13) Acenaphthene-d10	14.073	164	8353	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	18003	0.400	ng	# 0.00
29) Chrysene-d12	21.053	240	13959	0.400	ng	# 0.00
35) Perylene-d12	23.171	264	11697	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	2417	0.174	ng	0.00
5) Phenol-d6	6.629	99	1414	0.088	ng	0.00
8) Nitrobenzene-d5	8.579	82	2741	0.264	ng	0.00
11) 2-Methylnaphthalene-d10	11.788	152	5927	0.296	ng	0.00
14) 2,4,6-Tribromophenol	15.580	330	994	0.263	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	9439	0.278	ng	0.00
27) Fluoranthene-d10	18.876	212	18329	0.404	ng	-0.01
31) Terphenyl-d14	19.498	244	12125	0.435	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.068	88	12591	2.063	ng	# 86
6) bis(2-Chloroethyl)ether	6.874	93	673	0.047	ng	99

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

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