

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034141.D
 Acq On : 22 Sep 2024 01:47
 Operator : JU/RC
 Sample : P4116-26
 Misc :
 ALS Vial : 106 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE132D7-20240918

Quant Time: Sep 23 00:02:33 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	5620	0.400	ng	0.00
7) Naphthalene-d8	10.180	136	15180	0.400	ng	#-0.02
13) Acenaphthene-d10	14.073	164	8879	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	19085	0.400	ng	0.00
29) Chrysene-d12	21.044	240	12656	0.400	ng	#-0.01
35) Perylene-d12	23.169	264	9880	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	3106	0.195	ng	0.00
5) Phenol-d6	6.629	99	1900	0.102	ng	0.00
8) Nitrobenzene-d5	8.568	82	3419	0.294	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	7280	0.325	ng	0.00
14) 2,4,6-Tribromophenol	15.580	330	915	0.227	ng	0.00
15) 2-Fluorobiphenyl	12.683	172	11652	0.323	ng	-0.02
27) Fluoranthene-d10	18.876	212	18101	0.376	ng	-0.01
31) Terphenyl-d14	19.494	244	11681	0.462	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.068	88	19886	2.830	ng	# 84
6) bis(2-Chloroethyl)ether	6.875	93	421	0.026	ng	88

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

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