

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034142.D
 Acq On : 22 Sep 2024 02:23
 Operator : JU/RC
 Sample : P4116-27
 Misc :
 ALS Vial : 107 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP05-20240918

Quant Time: Sep 23 00:02:45 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	4792	0.400	ng	0.00
7) Naphthalene-d8	10.180	136	13672	0.400	ng	#-0.02
13) Acenaphthene-d10	14.072	164	6980	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	16412	0.400	ng	0.00
29) Chrysene-d12	21.044	240	12868	0.400	ng	#-0.01
35) Perylene-d12	23.168	264	3872	0.400	ng	#-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1577	0.116	ng	0.00
5) Phenol-d6	6.622	99	1303	0.082	ng	-0.01
8) Nitrobenzene-d5	8.568	82	2776	0.265	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	5854	0.290	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	818	0.259	ng	0.00
15) 2-Fluorobiphenyl	12.683	172	9445	0.333	ng	-0.02
27) Fluoranthene-d10	18.871	212	13663	0.330	ng	-0.02
31) Terphenyl-d14	19.494	244	9537	0.371	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.068	88	31900	5.325	ng	# 87
24) Pentachlorophenol	16.486	266	256	0.075	ng	# 91

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

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