

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120123\
 Data File : BN028963.D
 Acq On : 02 Dec 2023 00:49
 Operator : MA/JU
 Sample : 05358-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP03-20231108

Quant Time: Dec 02 03:16:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN113023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 23:11:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.797	152	2226	0.400	ng	0.00
7) Naphthalene-d8	10.584	136	6677	0.400	ng	0.00
13) Acenaphthene-d10	14.418	164	4106	0.400	ng	0.00
19) Phenanthrene-d10	17.171	188	9139	0.400	ng	# 0.00
29) Chrysene-d12	21.365	240	7502	0.400	ng	# 0.00
35) Perylene-d12	23.693	264	7240	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	643	0.095	ng	0.00
5) Phenol-d6	7.009	99	429	0.057	ng	0.00
8) Nitrobenzene-d5	8.961	82	1555	0.234	ng	-0.01
11) 2-Methylnaphthalene-d10	12.170	152	2515	0.235	ng	0.00
14) 2,4,6-Tribromophenol	15.918	330	757	0.273	ng	0.00
15) 2-Fluorobiphenyl	13.049	172	6649	0.375	ng	0.00
27) Fluoranthene-d10	19.204	212	8507	0.354	ng	0.00
31) Terphenyl-d14	19.808	244	8618	0.402	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.333	88	9736	4.466	ng	95
6) bis(2-Chloroethyl)ether	7.248	93	244	0.039	ng	88
16) Acenaphthylene	14.140	152	688	0.032	ng	98
30) Pyrene	19.599	202	302	0.099	ng	# 74

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

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