

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112124\
 Data File : BN035221.D
 Acq On : 21 Nov 2024 14:35
 Operator : RC/JU
 Sample : P4868-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW5-SP201-20241114

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/22/2024
 Supervised By :mohammad ahmed 11/22/2024

Quant Time: Nov 21 15:00:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN111324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 21 14:00:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.329	152	2998	0.400	ng	# 0.00
7) Naphthalene-d8	10.073	136	6728	0.400	ng	# 0.00
13) Acenaphthene-d10	13.983	164	1733	0.400	ng	0.00
19) Phenanthrene-d10	16.753	188	6277	0.400	ng	# 0.00
29) Chrysene-d12	21.006	240	144	0.400	ng	# 0.00
35) Perylene-d12	23.102	264	16	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.982	112	1218	0.160	ng	0.00
5) Phenol-d6	6.528	99	1253	0.131	ng	0.00
8) Nitrobenzene-d5	8.461	82	1475m	0.252	ng	0.00
11) 2-Methylnaphthalene-d10	11.678	152	3332	0.278	ng	0.00
14) 2,4,6-Tribromophenol	15.501	330	1252	1.001	ng	0.00
15) 2-Fluorobiphenyl	12.604	172	2481	0.353	ng	0.00
27) Fluoranthene-d10	18.808	212	2890	0.150	ng	0.00
31) Terphenyl-d14	19.435	244	4516	14.946	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.018	88	395	0.145	ng	# 49
24) Pentachlorophenol	16.406	266	86	0.044	ng	92

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

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