

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034496.D
 Acq On : 09 Oct 2024 08:33
 Operator : RC/JU
 Sample : P4226-21
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE117D1-20240926

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/10/2024
 Supervised By :mohammad ahmed 10/10/2024

Quant Time: Oct 09 08:57:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 26 14:44:02 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.380	152	3499	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	10070	0.400	ng	-0.01
13) Acenaphthene-d10	14.029	164	6140	0.400	ng	-0.02
19) Phenanthrene-d10	16.796	188	11730	0.400	ng	0.00
29) Chrysene-d12	21.008	240	7096	0.400	ng	# 0.00
35) Perylene-d12	23.121	264	7634	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.025	112	1413	0.142	ng	0.00
5) Phenol-d6	6.607	99	725m	0.063	ng	0.00
8) Nitrobenzene-d5	8.535	82	2082m	0.270	ng	0.00
11) 2-Methylnaphthalene-d10	11.742	152	5041	0.339	ng	-0.01
14) 2,4,6-Tribromophenol	15.542	330	887	0.319	ng	0.00
15) 2-Fluorobiphenyl	12.650	172	7306	0.293	ng	0.00
27) Fluoranthene-d10	18.834	212	12264	0.414	ng	-0.01
31) Terphenyl-d14	19.451	244	8436	0.595	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.032	88	1487	0.340	ng	# 60

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

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