

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034545.D
 Acq On : 11 Oct 2024 06:51
 Operator : RC/JU
 Sample : P4282-04
 Misc :
 ALS Vial : 121 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW7-SP303-20241002

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/11/2024
 Supervised By :mohammad ahmed 10/14/2024

Quant Time: Oct 11 07:18:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 00:21:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	3686	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	10574	0.400	ng	0.01
13) Acenaphthene-d10	14.026	164	6142	0.400	ng	0.01
19) Phenanthrene-d10	16.791	188	11579	0.400	ng	0.01
29) Chrysene-d12	21.006	240	7626	0.400	ng	0.00
35) Perylene-d12	23.114	264	9144	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1289	0.123	ng	0.00
5) Phenol-d6	6.636	99	658m	0.054	ng	0.07
8) Nitrobenzene-d5	8.547	82	2477m	0.306	ng	0.03
11) 2-Methylnaphthalene-d10	11.746	152	5493m	0.352	ng	0.02
14) 2,4,6-Tribromophenol	15.549	330	614	0.221	ng	0.01
15) 2-Fluorobiphenyl	12.647	172	7650	0.306	ng	0.01
27) Fluoranthene-d10	18.827	212	12706	0.435	ng	0.00
31) Terphenyl-d14	19.450	244	9144	0.600	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.025	88	443	0.096	ng	# 54

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

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