

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120924\
 Data File : BN035527.D
 Acq On : 10 Dec 2024 10:45
 Operator : RC/JU
 Sample : P5166-04DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW5-SP100-20241206DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/11/2024
 Supervised By :mohammad ahmed 12/11/2024

Quant Time: Dec 10 12:28:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN112724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 27 23:03:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.301	152	1780	0.400	ng	0.00
7) Naphthalene-d8	10.041	136	4601	0.400	ng	-0.01
13) Acenaphthene-d10	13.957	164	3512	0.400	ng	-0.01
19) Phenanthrene-d10	16.736	188	9370	0.400	ng	# 0.00
29) Chrysene-d12	20.974	240	6994	0.400	ng	0.00
35) Perylene-d12	23.067	264	6330	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.961	112	130	0.029	ng	0.00
5) Phenol-d6	6.506	99	82	0.015	ng	0.00
8) Nitrobenzene-d5	8.440	82	176m	0.063	ng	0.00
11) 2-Methylnaphthalene-d10	11.671	152	499	0.069	ng	0.02
14) 2,4,6-Tribromophenol	15.496	330	104	0.042	ng	0.02
15) 2-Fluorobiphenyl	12.589	172	825	0.062	ng	0.02
27) Fluoranthene-d10	18.785	212	2053	0.077	ng	0.00
31) Terphenyl-d14	19.412	244	1573	0.114	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.996	88	2988	1.756	ng	Qvalue 99

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

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