

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034549.D
 Acq On : 11 Oct 2024 09:15
 Operator : RC/JU
 Sample : P4263-04DL 2X
 Misc :
 ALS Vial : 125 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE115D2-20240930DL

Quant Time: Oct 11 10:59:38 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

Manual Integrations
 APPROVED

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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.372	152	4287	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	12310	0.400	ng	0.01
13) Acenaphthene-d10	14.029	164	7425	0.400	ng	0.01
19) Phenanthrene-d10	16.795	188	14129	0.400	ng	0.02
29) Chrysene-d12	21.008	240	8780	0.400	ng	0.00
35) Perylene-d12	23.112	264	10183	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	993	0.082	ng	0.00
5) Phenol-d6	6.643	99	517m	0.037	ng	0.07
8) Nitrobenzene-d5	8.557	82	1045m	0.111	ng	0.04
11) 2-Methylnaphthalene-d10	11.757	152	2574m	0.142	ng	0.03
14) 2,4,6-Tribromophenol	15.554	330	333	0.099	ng	0.02
15) 2-Fluorobiphenyl	12.650	172	3666	0.121	ng	0.01
27) Fluoranthene-d10	18.829	212	6432	0.180	ng	0.00
31) Terphenyl-d14	19.447	244	4414	0.252	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.024	88	17928	3.345	ng	Qvalue 99

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

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