

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111325\
 Data File : BN038200.D
 Acq On : 13 Nov 2025 13:27
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 50

Instrument :
 BNA_N
 ClientSampleId :
 SEEP-1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/13/2025
 Supervised By :Sohil Jodhani 11/14/2025

Quant Time: Nov 13 14:19:24 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 29 13:09:46 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.768	152	538910	20.000	ng	0.00
7) Naphthalene-d8	10.573	136	1785287	20.000	ng	0.00
13) Acenaphthene-d10	14.430	164	1021405	20.000	ng	0.00
19) Phenanthrene-d10	17.186	188	1654430	20.000	ng	# 0.00
29) Chrysene-d12	21.375	240	1150690	20.000	ng	# 0.00
35) Perylene-d12	23.680	264	1246822	20.000	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.341	112	4784	0.188	ng	0.00
5) Phenol-d6	6.937	99	8931m	0.289	ng	0.00
8) Nitrobenzene-d5	8.918	82	8631	0.300	ng	-0.01
11) 2-Methylnaphthalene-d10	12.167	152	14761	0.289	ng	-0.01
14) 2,4,6-Tribromophenol	15.932	330	2878	0.342	ng	0.00
15) 2-Fluorobiphenyl	13.051	172	21384	0.261	ng	0.00
27) Fluoranthene-d10	19.215	212	20805	0.249	ng	0.00
31) Terphenyl-d14	19.815	244	15374	0.307	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.247	88	125586	11.277	ng	Qvalue 97
24) Pentachlorophenol	16.838	266	301	0.028	ng	89
32) Benzo(a)anthracene	21.366	228	5577m	0.068	ng	
34) Bis(2-ethylhexyl)phtha...	21.286	149	14735	0.341	ng	# 89

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

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