

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035022.D
 Acq On : 12 Nov 2024 01:26
 Operator : RC/JU
 Sample : P4636-18
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CC0R2

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/12/2024
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 12 01:50:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.566	152	2897	0.400	ng/ul	0.00
4) Naphthalene-d8	10.328	136	7207	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.198	164	4940	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	10911	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.142	240	10522	0.400	ng/ul	# 0.00
23) Perylene-d12	23.308	264	11249m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	9007	3.267	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	2853	0.258	ng/ul	-0.01
18) Fluoranthene-d10	18.981	212	7649	0.283	ng/ul	0.00
Target Compounds						
14) Pentachlorophenol	16.603	266	126	0.050	ng/ul#	86
19) Fluoranthene	19.009	202	1264	0.036	ng/ul#	74
20) Pyrene	19.376	202	1312	0.036	ng/ul#	75
21) Benzo(a)anthracene	21.128	228	848	0.022	ng/ul#	62
22) Chrysene	21.180	228	946	0.024	ng/ul#	64
24) Benzo(b)fluoranthene	22.665	252	1205m	0.030	ng/ul	
25) Benzo(k)fluoranthene	22.694	252	1150m	0.028	ng/ul	

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

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