

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035012.D
 Acq On : 11 Nov 2024 19:26
 Operator : RC/JU
 Sample : P4636-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CC0P2

Manual Integrations
 APPROVED

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

Quant Time: Nov 11 22:04:18 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	7.562	152	2894	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.328	136	7681	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.198	164	5527	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	12573	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	12422	0.400	ng/ul	# 0.00
23) Perylene-d12	23.308	264	12535m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	8097	2.940	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	2684	0.228	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	7911	0.248	ng/ul	0.00
Target Compounds						
15) Phenanthrene	16.987	178	1298	0.039	ng/ul#	90
19) Fluoranthene	19.009	202	3501	0.084	ng/ul#	89
20) Pyrene	19.372	202	3396m	0.079	ng/ul	
21) Benzo(a)anthracene	21.125	228	2839	0.063	ng/ul#	89
22) Chrysene	21.177	228	3317	0.072	ng/ul#	90
24) Benzo(b)fluoranthene	22.665	252	5336m	0.121	ng/ul	
25) Benzo(k)fluoranthene	22.700	252	2015m	0.044	ng/ul	
26) Benzo(a)pyrene	23.214	252	3236m	0.082	ng/ul	
27) Indeno(1,2,3-cd)pyrene	25.453	276	2602m	0.054	ng/ul	
29) Benzo(g,h,i)perylene	26.107	276	2689m	0.073	ng/ul	

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

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