

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035013.D
 Acq On : 11 Nov 2024 20:02
 Operator : RC/JU
 Sample : P4636-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CC0P4

Manual Integrations
 APPROVED

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

Quant Time: Nov 11 22:04:27 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	2914	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.328	136	7635	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.199	164	5369	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	11943	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	11581	0.400	ng/ul	# 0.00
23) Perylene-d12	23.308	264	12033m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	8430	3.040	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	2599	0.222	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	7480	0.252	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.603	266	71	0.026	ng/ul#	84
15) Phenanthrene	16.987	178	3151	0.099	ng/ul	95
16) Anthracene	17.080	178	1014	0.034	ng/ul#	80
19) Fluoranthene	19.009	202	3130	0.081	ng/ul#	89
20) Pyrene	19.372	202	2317	0.058	ng/ul#	82
21) Benzo(a)anthracene	21.125	228	871	0.021	ng/ul#	68
22) Chrysene	21.177	228	1531m	0.036	ng/ul	
24) Benzo(b)fluoranthene	22.665	252	1103m	0.026	ng/ul	
29) Benzo(g,h,i)perylene	26.110	276	785m	0.022	ng/ul	

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

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