

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103124\
 Data File : BN034780.D
 Acq On : 31 Oct 2024 23:42
 Operator : RC/JU
 Sample : P4529-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F7H15

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024

Quant Time: Nov 03 23:23:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 11:13:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.600	152	7471	0.400	ng/ul	0.00
4) Naphthalene-d8	10.366	136	23626	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.236	164	13350	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.979	188	26631	0.400	ng/ul	0.00
17) Chrysene-d12	21.172	240	18341	0.400	ng/ul	0.00
23) Perylene-d12	23.355	264	15596m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.170	96	25739	3.328	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.961	152	6372	0.215	ng/ul	0.00
18) Fluoranthene-d10	19.014	212	12346	0.207	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Anthracene	17.110	178	3157	0.042	ng/ul#	92
19) Fluoranthene	19.042	202	8997	0.101	ng/ul	96
20) Pyrene	19.404	202	18221m	0.197	ng/ul	
21) Benzo(a)anthracene	21.157	228	3500	0.048	ng/ul	93
22) Chrysene	21.207	228	5399	0.073	ng/ul	96
24) Benzo(b)fluoranthene	22.706	252	7339m	0.116	ng/ul	
25) Benzo(k)fluoranthene	22.744	252	2561m	0.040	ng/ul	
26) Benzo(a)pyrene	23.264	252	2661	0.049	ng/ul#	44
27) Indeno(1,2,3-cd)pyrene	25.527	276	1277	0.022	ng/ul	99
29) Benzo(g,h,i)perylene	26.188	276	996	0.022	ng/ul#	57

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

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