

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035160.D
 Acq On : 19 Nov 2024 13:53
 Operator : RC/JU
 Sample : P4847-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A0AT1

Quant Time: Nov 19 14:21:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.545	152	3059	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	7349	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.185	164	5295	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.941	188	11358	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	8770	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	8504	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1564	0.559	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	4049	0.376	ng/ul	0.00
18) Fluoranthene-d10	18.977	212	11096	0.459	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.355	128	42740	2.309	ng/ul	97
10) Acenaphthylene	13.903	152	291917	13.898	ng/ul	98
11) Acenaphthene	14.249	153	200527	12.478	ng/ul	98
12) Fluorene	15.244	166	43311	2.296	ng/ul	99
15) Phenanthrene	16.979	178	128847	4.291	ng/ul	99
16) Anthracene	17.072	178	125158	4.491	ng/ul	98
19) Fluoranthene	19.009	202	126930	4.157	ng/ul	98
20) Pyrene	19.372	202	128507	4.044	ng/ul	100
21) Benzo(a)anthracene	21.134	228	56693	1.863	ng/ul	100
22) Chrysene	21.183	228	39605	1.293	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	130940	4.658	ng/ul	91
25) Benzo(k)fluoranthene	22.717	252	58918	2.033	ng/ul#	91
26) Benzo(a)pyrene	23.226	252	47068	1.832	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.480	276	46662	1.345	ng/ul	97
28) Dibenzo(a,h)anthracene	25.493	278	47745	1.755	ng/ul	93
29) Benzo(g,h,i)perylene	26.131	276	41242	1.508	ng/ul	97

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

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