

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070124\
 Data File : BN032358.D
 Acq On : 01 Jul 2024 12:12
 Operator : MA/JU
 Sample : P2934-25
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN5

Quant Time: Jul 01 12:37:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 13:23:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.612	152	6755	0.400	ng/ul	0.00
4) Naphthalene-d8	10.387	136	20807	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.258	164	11020	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.012	188	19576	0.400	ng/ul	-0.01
17) Chrysene-d12	21.214	240	15461	0.400	ng/ul	#-0.01
23) Perylene-d12	23.433	264	18258	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	3375	0.449	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.988	152	9189	0.294	ng/ul	-0.01
18) Fluoranthene-d10	19.045	212	16497	0.355	ng/ul	-0.01
Target Compounds						Qvalue
5) Naphthalene	10.431	128	971103	17.636	ng/ul	99
10) Acenaphthylene	13.975	152	854667	18.648	ng/ul	99
11) Acenaphthene	14.318	153	518616	14.580	ng/ul	98
12) Fluorene	15.312	166	309954	7.313	ng/ul	98
15) Phenanthrene	17.050	178	436511	8.133	ng/ul	99
16) Anthracene	17.143	178	311710	6.511	ng/ul	99
19) Fluoranthene	19.078	202	533822	9.067	ng/ul	99
20) Pyrene	19.445	202	644705	10.479	ng/ul	97
21) Benzo(a)anthracene	21.197	228	182210	3.265	ng/ul	100
22) Chrysene	21.249	228	324604	5.300	ng/ul	99
24) Benzo(b)fluoranthene	22.763	252	402479	5.074	ng/ul	93
25) Benzo(k)fluoranthene	22.807	252	335305	4.351	ng/ul	95
26) Benzo(a)pyrene	23.336	252	436736	7.913	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.663	276	538785	6.235	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.676	278	343495	5.350	ng/ul	96
29) Benzo(g,h,i)perylene	26.354	276	378692	5.501	ng/ul	98

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

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