

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062724\
 Data File : BN032282.D
 Acq On : 27 Jun 2024 16:13
 Operator : MA/JU
 Sample : P2909-16DL 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN6DL

Quant Time: Jun 27 16:40:13 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	7.620	152	6782	0.400	ng/ul	0.00
4) Naphthalene-d8	10.398	136	21725	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.267	164	15735	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.020	188	35591	0.400	ng/ul	0.00
17) Chrysene-d12	21.223	240	29637	0.400	ng/ul	0.00
23) Perylene-d12	23.447	264	27213	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.156	96	489	0.065	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.999	152	1102	0.034	ng/ul	0.00
18) Fluoranthene-d10	19.055	212	3463	0.039	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.448	128	29087	0.506	ng/ul	99
10) Acenaphthylene	13.985	152	50109	0.766	ng/ul	99
11) Acenaphthene	14.327	153	85014	1.674	ng/ul	98
12) Fluorene	15.322	166	17130	0.283	ng/ul	100
15) Phenanthrene	17.058	178	170200	1.744	ng/ul	99
16) Anthracene	17.155	178	133472	1.534	ng/ul	99
19) Fluoranthene	19.087	202	98723	0.875	ng/ul	99
20) Pyrene	19.450	202	122069	1.035	ng/ul	99
21) Benzo(a)anthracene	21.209	228	45004	0.421	ng/ul	99
22) Chrysene	21.261	228	82267	0.701	ng/ul	99
24) Benzo(b)fluoranthene	22.778	252	95040	0.804	ng/ul	96
25) Benzo(k)fluoranthene	22.822	252	66229	0.577	ng/ul	98
26) Benzo(a)pyrene	23.351	252	9967	0.121	ng/ul#	65
27) Indeno(1,2,3-cd)pyrene	25.683	276	69827	0.542	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.696	278	55642	0.581	ng/ul	98
29) Benzo(g,h,i)perylene	26.371	276	48749	0.475	ng/ul	99

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

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