

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071624\
 Data File : BN032738.D
 Acq On : 16 Jul 2024 13:25
 Operator : MA/JU
 Sample : P3177-22DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DQ2DL

Quant Time: Jul 16 13:50:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.561	152	3534	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.342	136	10696	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.201	164	5944	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.964	188	12070	0.400	ng/ul	-0.03
17) Chrysene-d12	21.184	240	10157	0.400	ng/ul	-0.01
23) Perylene-d12	23.379	264	12991	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.105	96	179	0.042	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.964	152	312	0.019	ng/ul	0.00
18) Fluoranthene-d10	19.002	212	1004	0.030	ng/ul	-0.02
Target Compounds						
					Qvalue	
5) Naphthalene	10.392	128	17369	0.605	ng/ul	99
10) Acenaphthylene	13.923	152	17643	0.735	ng/ul	100
11) Acenaphthene	14.261	153	23747	1.196	ng/ul	99
12) Fluorene	15.274	166	13431	0.592	ng/ul	100
15) Phenanthrene	17.006	178	43836	1.334	ng/ul	99
16) Anthracene	17.103	178	38455	1.432	ng/ul	98
19) Fluoranthene	19.035	202	33563	0.813	ng/ul	99
20) Pyrene	19.402	202	14936	0.349	ng/ul	100
21) Benzo(a)anthracene	21.172	228	3867	0.120	ng/ul	93
22) Chrysene	21.219	228	14313	0.306	ng/ul	97
24) Benzo(b)fluoranthene	22.721	252	13911	0.303	ng/ul	88
25) Benzo(k)fluoranthene	22.765	252	13176	0.236	ng/ul#	83
26) Benzo(a)pyrene	23.288	252	21070	0.531	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.595	276	24007	0.447	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.598	278	22714	0.546	ng/ul	98
29) Benzo(g,h,i)perylene	26.269	276	16514	0.368	ng/ul	98

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

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