

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020625\
 Data File : BN036349.D
 Acq On : 06 Feb 2025 23:22
 Operator : RC/JU
 Sample : PB166543BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS543

Quant Time: Feb 07 07:50:52 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN012125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 14:44:03 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.771	152	2119	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.562	136	4738	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.409	164	2517	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.154	188	4928	0.400	ng/ul	#-0.02
17) Chrysene-d12	21.342	240	4327	0.400	ng/ul	#-0.02
23) Perylene-d12	23.622	264	4632	0.400	ng/ul	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	1731	0.746	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.157	152	2327	0.389	ng/ul	0.00
18) Fluoranthene-d10	19.184	212	4483	0.377	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.286	88	4918	1.986	ng/ul	88
5) Naphthalene	10.612	128	5162	0.395	ng/ul	99
7) 2-Methylnaphthalene	12.229	142	3158	0.389	ng/ul	100
8) 1-Methylnaphthalene	12.449	142	3219	0.391	ng/ul	100
10) Acenaphthylene	14.127	152	4214	0.355	ng/ul	99
11) Acenaphthene	14.474	153	3357	0.384	ng/ul	98
12) Fluorene	15.464	166	3533	0.366	ng/ul	97
14) Pentachlorophenol	16.817	266	1098	0.636	ng/ul	99
15) Phenanthrene	17.196	178	5487	0.380	ng/ul	99
16) Anthracene	17.289	178	4742	0.354	ng/ul	98
19) Fluoranthene	19.216	202	6170	0.375	ng/ul#	95
20) Pyrene	19.579	202	6451	0.376	ng/ul#	93
21) Benzo(a)anthracene	21.324	228	5695	0.369	ng/ul	100
22) Chrysene	21.377	228	6233	0.395	ng/ul	99
24) Benzo(b)fluoranthene	22.935	252	6149	0.398	ng/ul	86
25) Benzo(k)fluoranthene	22.978	252	6382	0.393	ng/ul#	85
26) Benzo(a)pyrene	23.519	252	5724	0.410	ng/ul#	79
27) Indeno(1,2,3-cd)pyrene	25.940	276	7265	0.393	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.954	278	5565	0.391	ng/ul#	90
29) Benzo(g,h,i)perylene	26.648	276	5959	0.416	ng/ul	96

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

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