

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035574.D
 Acq On : 11 Dec 2024 19:42
 Operator : RC/JU
 Sample : PB165482BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS482

Quant Time: Dec 11 22:10:20 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 : Tue Dec 10 16:10:57 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.293	152	2520	0.400	ng/ul	0.00
4) Naphthalene-d8	10.036	136	7352	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	4997	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	10510	0.400	ng/ul	0.00
17) Chrysene-d12	20.964	240	7880	0.400	ng/ul	0.00
23) Perylene-d12	23.054	264	7629	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	1441	0.626	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	3760	0.349	ng/ul	0.00
18) Fluoranthene-d10	18.767	212	8166	0.376	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	3963	1.585	ng/ul	93
5) Naphthalene	10.086	128	6354	0.343	ng/ul	99
7) 2-Methylnaphthalene	11.714	142	4037	0.314	ng/ul	97
8) 1-Methylnaphthalene	11.939	142	4195	0.324	ng/ul	98
10) Acenaphthylene	13.662	152	6296	0.318	ng/ul	99
11) Acenaphthene	14.013	153	4885	0.322	ng/ul	99
12) Fluorene	15.017	166	5457	0.307	ng/ul	99
14) Pentachlorophenol	16.379	266	1706	0.697	ng/ul	99
15) Phenanthrene	16.759	178	8852	0.319	ng/ul	100
16) Anthracene	16.852	178	7939	0.308	ng/ul	100
19) Fluoranthene	18.800	202	9598	0.350	ng/ul	97
20) Pyrene	19.167	202	10114	0.354	ng/ul#	96
21) Benzo(a)anthracene	20.947	228	8806	0.322	ng/ul	99
22) Chrysene	20.999	228	9435	0.343	ng/ul	100
24) Benzo(b)fluoranthene	22.440	252	8423	0.334	ng/ul	98
25) Benzo(k)fluoranthene	22.483	252	9060	0.348	ng/ul	98
26) Benzo(a)pyrene	22.966	252	7646	0.332	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.084	276	8855	0.285	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.104	278	6771	0.277	ng/ul#	95
29) Benzo(g,h,i)perylene	25.698	276	7039	0.287	ng/ul#	96

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

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