

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121724\
 Data File : BM049025.D
 Acq On : 17 Dec 2024 17:02
 Operator : RC/JU
 Sample : PB165589BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS589

Quant Time: Dec 18 01:17:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM120924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 17 10:03:27 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	9239	0.400	ng/ul	0.00
4) Naphthalene-d8	10.336	136	28242	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.205	164	15790	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	29106	0.400	ng/ul #	0.00
17) Chrysene-d12	21.149	240	19851	0.400	ng/ul #	0.00
23) Perylene-d12	23.306	264	20681	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.155	96	5728	0.530	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.931	152	12284	0.341	ng/ul	0.00
18) Fluoranthene-d10	18.987	212	20451	0.345	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.185	88	16510	1.321	ng/ul#	81
5) Naphthalene	10.385	128	21054	0.291	ng/ul	100
7) 2-Methylnaphthalene	12.008	142	13277	0.300	ng/ul	99
8) 1-Methylnaphthalene	12.228	142	13667	0.303	ng/ul	100
10) Acenaphthylene	13.922	152	18887	0.283	ng/ul	98
11) Acenaphthene	14.265	153	14194	0.281	ng/ul	98
12) Fluorene	15.259	166	15336	0.274	ng/ul	99
14) Pentachlorophenol	16.608	266	3364	0.561	ng/ul	94
15) Phenanthrene	16.992	178	22609	0.283	ng/ul	100
16) Anthracene	17.085	178	20738	0.292	ng/ul	99
19) Fluoranthene	19.015	202	25046	0.304	ng/ul	95
20) Pyrene	19.377	202	26014	0.293	ng/ul	96
21) Benzo(a)anthracene	21.134	228	21159	0.303	ng/ul	98
22) Chrysene	21.184	228	23466	0.297	ng/ul	100
24) Benzo(b)fluoranthene	22.666	252	21127	0.296	ng/ul	97
25) Benzo(k)fluoranthene	22.709	252	22982	0.278	ng/ul#	94
26) Benzo(a)pyrene	23.212	252	19769	0.299	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.447	276	26292	0.283	ng/ul#	48
28) Dibenzo(a,h)anthracene	25.467	278	20427	0.283	ng/ul	95
29) Benzo(g,h,i)perylene	26.098	276	21158	0.283	ng/ul	96

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

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