

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103124\
 Data File : BM048413.D
 Acq On : 02 Nov 2024 02:34
 Operator : RC/JU
 Sample : PB164517BS
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS517

Quant Time: Nov 03 21:51:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	2982	0.400	ng/ul	0.04
4) Naphthalene-d8	10.546	136	11575	0.400	ng/ul	0.02
9) Acenaphthene-d10	14.368	164	6879	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.129	188	14538	0.400	ng/ul	0.00
17) Chrysene-d12	21.298	240	12484	0.400	ng/ul	0.00
23) Perylene-d12	23.534	264	12770	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	3314	0.905	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.168	152	5148	0.350	ng/ul	0.03
18) Fluoranthene-d10	19.146	212	12936	0.408	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.198	88	9007	2.267	ng/ul#	83
5) Naphthalene	10.601	128	9661	0.319	ng/ul	97
7) 2-Methylnaphthalene	12.240	142	5371	0.298	ng/ul	97
8) 1-Methylnaphthalene	12.432	142	6473	0.337	ng/ul	98
10) Acenaphthylene	14.095	152	8485	0.319	ng/ul	98
11) Acenaphthene	14.433	153	7117	0.334	ng/ul	98
12) Fluorene	15.446	166	7673	0.324	ng/ul	98
14) Pentachlorophenol	16.795	266	2381	0.588	ng/ul	94
15) Phenanthrene	17.171	178	11378	0.315	ng/ul	99
16) Anthracene	17.289	178	10156	0.300	ng/ul	98
19) Fluoranthene	19.174	202	15906	0.370	ng/ul	98
20) Pyrene	19.537	202	17071	0.370	ng/ul	99
21) Benzo(a)anthracene	21.286	228	11686	0.312	ng/ul	100
22) Chrysene	21.330	228	19514	0.379	ng/ul	99
24) Benzo(b)fluoranthene	22.858	252	14626	0.335	ng/ul	100
25) Benzo(k)fluoranthene	22.905	252	19323	0.374	ng/ul	100
26) Benzo(a)pyrene	23.446	252	14372	0.351	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.833	276	18851	0.329	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.856	278	14553	0.326	ng/ul	97
29) Benzo(g,h,i)perylene	26.507	276	16420	0.340	ng/ul	99

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

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