

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021225\
 Data File : BM049638.D
 Acq On : 13 Feb 2025 02:55
 Operator : RC/JU
 Sample : PB166627BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS627

Quant Time: Feb 17 12:06:17 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM021225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 12 16:40:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.813	152	11011	0.400	ng/ul	0.00
4) Naphthalene-d8	10.607	136	33442	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.451	164	20846	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.188	188	43892	0.400	ng/ul	0.00
17) Chrysene-d12	21.365	240	32502	0.400	ng/ul	0.00
23) Perylene-d12	23.650	264	35732	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	8212	0.663	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.196	152	20856	0.454	ng/ul	0.00
18) Fluoranthene-d10	19.216	212	43926	0.444	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	22266	1.663	ng/ul	86
5) Naphthalene	10.656	128	32981	0.398	ng/ul	100
7) 2-Methylnaphthalene	12.273	142	22276	0.413	ng/ul	99
8) 1-Methylnaphthalene	12.493	142	22938	0.419	ng/ul	99
10) Acenaphthylene	14.169	152	35022	0.396	ng/ul	99
11) Acenaphthene	14.511	153	24810	0.385	ng/ul	98
12) Fluorene	15.497	166	29057	0.394	ng/ul	98
14) Pentachlorophenol	16.841	266	10339	1.149	ng/ul	97
15) Phenanthrene	17.230	178	46514	0.396	ng/ul	100
16) Anthracene	17.323	178	44400	0.413	ng/ul	99
19) Fluoranthene	19.244	202	53137	0.396	ng/ul	99
20) Pyrene	19.606	202	56469	0.408	ng/ul	99
21) Benzo(a)anthracene	21.347	228	47982	0.431	ng/ul	99
22) Chrysene	21.400	228	47047	0.398	ng/ul	99
24) Benzo(b)fluoranthene	22.960	252	48234	0.405	ng/ul	97
25) Benzo(k)fluoranthene	23.007	252	50719	0.379	ng/ul	98
26) Benzo(a)pyrene	23.551	252	44351	0.413	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.977	276	59106	0.388	ng/ul	99
28) Dibenzo(a,h)anthracene	25.990	278	45692	0.395	ng/ul	98
29) Benzo(g,h,i)perylene	26.684	276	46232	0.378	ng/ul	98

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

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