

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110723\
 Data File : BM042642.D
 Acq On : 08 Nov 2023 09:45
 Operator : MA/JU
 Sample : 05168-03
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 XA014

Quant Time: Nov 08 10:17:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 02:23:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.868	152	1090	0.400	ng/ul	0.00
4) Naphthalene-d8	10.671	136	2682	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.510	164	1603	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.262	188	3526	0.400	ng/ul	0.00
17) Chrysene-d12	21.444	240	2898	0.400	ng/ul	#-0.01
23) Perylene-d12	23.829	264	3005	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	962	0.668	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.277	152	1325	0.354	ng/ul	0.00
18) Fluoranthene-d10	19.289	212	3811	0.380	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.721	128	59164	6.832	ng/ul#	93
10) Acenaphthylene	14.237	152	23703	2.161	ng/ul	97
11) Acenaphthene	14.570	153	78475	10.543	ng/ul	98
12) Fluorene	15.560	166	37858	4.434	ng/ul	99
15) Phenanthrene	17.304	178	31760	2.323	ng/ul	96
16) Anthracene	17.397	178	39482	3.191	ng/ul	95
19) Fluoranthene	19.317	202	60057	3.316	ng/ul	97
20) Pyrene	19.684	202	20167	1.013	ng/ul	97
21) Benzo(a)anthracene	21.426	228	48401	3.752	ng/ul	96
22) Chrysene	21.479	228	36913	2.768	ng/ul	97
24) Benzo(b)fluoranthene	23.086	252	51981	4.283	ng/ul	79
25) Benzo(k)fluoranthene	23.136	252	38279	3.010	ng/ul#	79
26) Benzo(a)pyrene	23.718	252	43984	3.500	ng/ul#	72
27) Indeno(1,2,3-cd)pyrene	26.309	276	23132	1.416	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.319	278	32181	2.759	ng/ul#	75
29) Benzo(g,h,i)perylene	27.077	276	15462	1.018	ng/ul#	90

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

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