

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071323\
 Data File : BM040854.D
 Acq On : 13 Jul 2023 17:59
 Operator : MA/JU
 Sample : 03418-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4663

Quant Time: Jul 14 03:16:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 14 03:14:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.866	152	4946	0.400	ng/ul	0.00
4) Naphthalene-d8	10.669	136	17823	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.513	164	8001	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.265	188	15172	0.400	ng/ul	0.00
17) Chrysene-d12	21.457	240	9552	0.400	ng/ul	# 0.00
23) Perylene-d12	23.842	264	8602	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	27870	3.587	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.264	152	6475	0.259	ng/ul	0.00
18) Fluoranthene-d10	19.297	212	11093	0.362	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.719	128	730183	14.853	ng/ul	99
8) 1-Methylnaphthalene	12.550	142	622843	21.586	ng/ul	100
11) Acenaphthene	14.573	153	621737	20.598	ng/ul	98
12) Fluorene	15.564	166	9025	0.280	ng/ul#	96
14) Pentachlorophenol	16.915	266	268373	91.866	ng/ul	93
15) Phenanthrene	17.308	178	888043	18.843	ng/ul	98
16) Anthracene	17.396	178	1016449	25.666	ng/ul	99
19) Fluoranthene	19.329	202	2444101	60.237	ng/ul#	91
20) Pyrene	19.692	202	1547616	35.577	ng/ul	95
21) Benzo(a)anthracene	21.442	228	1689530	52.292	ng/ul	99
22) Chrysene	21.495	228	964420	26.167	ng/ul	99
24) Benzo(b)fluoranthene	23.114	252	1019676	30.630	ng/ul	85
26) Benzo(a)pyrene	23.737	252	835967	27.878	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.307	276	1200150	28.738	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.320	278	754821	23.085	ng/ul#	91

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

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