

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072623\
 Data File : BM041129.D
 Acq On : 26 Jul 2023 11:06
 Operator : MA/JU
 Sample : 03553-10MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4701MS

Quant Time: Jul 27 01:17:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM072423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 04:20:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.836	152	2672	0.400	ng/ul	0.00
4) Naphthalene-d8	10.641	136	9299	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.486	164	4821	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.248	188	9519	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	8791	0.400	ng/ul	0.00
23) Perylene-d12	23.833	264	8616	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.254	96	980	0.238	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.242	152	4522	0.341	ng/ul	0.00
18) Fluoranthene-d10	19.282	212	11057	0.462	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.288	88	2558	0.607	ng/ul#	54
5) Naphthalene	10.685	128	8710	0.288	ng/ul	99
7) 2-Methylnaphthalene	12.313	142	4905	0.331	ng/ul	99
8) 1-Methylnaphthalene	12.528	142	5386	0.354	ng/ul	99
10) Acenaphthylene	14.213	152	6838	0.336	ng/ul	100
11) Acenaphthene	14.550	153	5798	0.337	ng/ul	98
12) Fluorene	15.550	166	6365	0.339	ng/ul	99
14) Pentachlorophenol	16.919	266	384	0.139	ng/ul	93
15) Phenanthrene	17.291	178	10539	0.359	ng/ul	98
16) Anthracene	17.388	178	8553	0.360	ng/ul	99
19) Fluoranthene	19.315	202	13521	0.429	ng/ul	98
20) Pyrene	19.677	202	14332	0.413	ng/ul	99
21) Benzo(a)anthracene	21.434	228	11765	0.400	ng/ul	99
22) Chrysene	21.483	228	12946	0.381	ng/ul	100
24) Benzo(b)fluoranthene	23.105	252	13285	0.387	ng/ul	100
25) Benzo(k)fluoranthene	23.152	252	12661	0.386	ng/ul	100
26) Benzo(a)pyrene	23.728	252	10865	0.391	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.297	276	15602	0.379	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.314	278	12292	0.381	ng/ul	99
29) Benzo(g,h,i)perylene	27.051	276	13841	0.376	ng/ul	99

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

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