

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112723\
 Data File : BM043046.D
 Acq On : 27 Nov 2023 12:22
 Operator : MA/JU
 Sample : 05261-02DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCKV2DL

Quant Time: Nov 27 12:54:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 27 02:36:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.889	152	765	0.400	ng/ul	0.00
4) Naphthalene-d8	10.655	136	2623	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.454	164	1848	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.220	188	3978	0.400	ng/ul	-0.02
17) Chrysene-d12	21.415	240	2476	0.400	ng/ul	0.00
23) Perylene-d12	23.774	264	2609	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	136	0.116	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.283	152	70	0.020	ng/ul	0.03
18) Fluoranthene-d10	19.247	212	318	0.035	ng/ul	-0.01
Target Compounds						
					Qvalue	
5) Naphthalene	10.704	128	2237	0.314	ng/ul	92
10) Acenaphthylene	14.191	152	4643	0.492	ng/ul	97
11) Acenaphthene	14.519	153	7676	1.074	ng/ul	97
12) Fluorene	15.546	166	1341	0.183	ng/ul#	95
15) Phenanthrene	17.258	178	13246	1.167	ng/ul	93
16) Anthracene	17.359	178	10517	0.978	ng/ul	91
19) Fluoranthene	19.280	202	8552	0.609	ng/ul#	94
20) Pyrene	19.647	202	10576	0.708	ng/ul#	94
21) Benzo(a)anthracene	21.400	228	2575	0.321	ng/ul	98
22) Chrysene	21.450	228	7096	0.474	ng/ul	97
24) Benzo(b)fluoranthene	23.040	252	6684	0.759	ng/ul	88
25) Benzo(k)fluoranthene	23.095	252	6020	0.496	ng/ul	93
26) Benzo(a)pyrene	23.683	252	751	0.078	ng/ul#	52
27) Indeno(1,2,3-cd)pyrene	26.225	276	5956	0.458	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.239	278	4622	0.459	ng/ul#	89
29) Benzo(g,h,i)perylene	26.993	276	4745	0.375	ng/ul	97

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

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