

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027927.D
 Acq On : 28 Sep 2023 15:52
 Operator : MA/JU
 Sample : 04480-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Sep 28 17:18:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 27 03:52:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	3601	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	14161	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	9425	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	20802	0.400	ng/ul	0.00
17) Chrysene-d12	21.539	240	18852	0.400	ng/ul	# 0.00
23) Perylene-d12	24.012	264	18942	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	29342	6.469	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	11283	0.526	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	32732	0.568	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.368	88	1236	0.262	ng/ul#	82
7) 2-Methylnaphthalene	12.434	142	875	0.034	ng/ul	98
8) 1-Methylnaphthalene	12.654	142	684	0.026	ng/ul	96
14) Pentachlorophenol	16.991	266	494	0.101	ng/ul	95
15) Phenanthrene	17.396	178	2713	0.042	ng/ul#	90
16) Anthracene	17.489	178	2389	0.044	ng/ul#	88
19) Fluoranthene	19.413	202	3810	0.050	ng/ul#	92
20) Pyrene	19.780	202	3646	0.046	ng/ul#	91
21) Benzo(a)anthracene	21.522	228	4031	0.055	ng/ul#	74
22) Chrysene	21.577	228	3940	0.052	ng/ul#	75
24) Benzo(b)fluoranthene	23.246	252	3913	0.046	ng/ul#	1
25) Benzo(k)fluoranthene	23.293	252	3550	0.043	ng/ul#	1
26) Benzo(a)pyrene	23.904	252	2821	0.042	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.616	276	3966	0.046	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.636	278	3560	0.051	ng/ul#	1
29) Benzo(g,h,i)perylene	27.424	276	3304	0.045	ng/ul#	19

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

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