

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012993.D
 Acq On : 08 Dec 2022 19:57
 Operator : CG/JU
 Sample : N5894-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 09 01:58:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 08 00:35:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	568543	20.000	ng/ul	0.00
20) Naphthalene-d8	10.792	136	2477812	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	1682631	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	3850956	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3516022	20.000	ng/ul	0.00
88) Perylene-d12	23.945	264	3668594	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	73306	5.513	ng/uL	0.00
4) Pyridine-d5	3.787	84	334584	8.858	ng/ul	0.00
7) Phenol-d5	7.128	99	367271	7.931	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	1021404	36.563	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	1032757	28.330	ng/ul	0.00
15) 4-Methylphenol-d8	8.681	113	696112	18.346	ng/ul	0.00
21) Nitrobenzene-d5	9.145	128	671513	36.886	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	719398	35.100	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	1246588	31.314	ng/ul	0.00
31) 4-Chloroaniline-d4	10.928	131	1226017	21.815	ng/ul	0.00
46) Dimethylphthalate-d6	14.022	166	4788962	37.033	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	5212661	36.683	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	125639	5.401	ng/ul	-0.01
60) Fluorene-d10	15.610	176	4212693	38.506	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	749646	30.250	ng/ul	0.00
73) Anthracene-d10	17.468	188	6941251	39.953	ng/ul	0.00
81) Pyrene-d10	19.698	212	8280906	41.031	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.786	264	7429689	40.631	ng/ul	0.00

Target Compounds	Qvalue

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

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