

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012989.D
 Acq On : 08 Dec 2022 17:40
 Operator : CG/JU
 Sample : N5894-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA0

Quant Time: Dec 09 01:57:05 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	612225	20.000	ng/ul	0.00
20) Naphthalene-d8	10.793	136	2678799	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	1831017	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	4206624	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3858358	20.000	ng/ul	0.00
88) Perylene-d12	23.945	264	4089556	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	67002	4.679	ng/uL	0.00
4) Pyridine-d5	3.787	84	218477	5.371	ng/ul	0.00
7) Phenol-d5	7.128	99	243290	4.879	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	785983	26.128	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	737400	18.784	ng/ul	0.00
15) 4-Methylphenol-d8	8.681	113	477944	11.697	ng/ul	0.00
21) Nitrobenzene-d5	9.146	128	504857	25.651	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	509947	23.014	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	894049	20.773	ng/ul	0.00
31) 4-Chloroaniline-d4	10.928	131	953309	15.690	ng/ul	0.00
46) Dimethylphthalate-d6	14.022	166	4188625	29.765	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	4268546	27.605	ng/ul	0.00
54) 4-Nitrophenol-d4	14.804	143	93255	3.684	ng/ul	0.00
60) Fluorene-d10	15.610	176	3754408	31.536	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	654409	24.174	ng/ul	0.00
73) Anthracene-d10	17.469	188	6658193	35.084	ng/ul	0.00
81) Pyrene-d10	19.698	212	7958927	35.937	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.786	264	7297859	35.802	ng/ul	0.00

Target Compounds Qvalue

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

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