

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021225\
 Data File : BP024051.D
 Acq On : 12 Feb 2025 15:42
 Operator : RC/JU
 Sample : PB165831BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK831

Quant Time: Feb 14 11:22:24 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 13 12:56:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	510544	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	1926797	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.381	164	1078300	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	2083050	20.000	ng/u1	0.01
79) Chrysene-d12	21.604	240	1888352	20.000	ng/u1	0.00
88) Perylene-d12	24.962	264	1878400	20.000	ng/u1	0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	263	0.021	ng/uL	-0.11
4) Pyridine-d5	3.705	84	71	0.002	ng/u1	0.00
7) Phenol-d5	6.940	99	18	0.000	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.087	67	91	0.004	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.287	132	20	0.001	ng/u1	-0.02
15) 4-Methylphenol-d8	8.457	113	24	0.001	ng/u1	0.00
21) Nitrobenzene-d5	8.922	128	66	0.004	ng/u1	0.02
24) 2-Nitrophenol-d4	9.640	143	24	0.001	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.157	165	41	0.001	ng/u1	0.00
31) 4-Chloroaniline-d4	10.687	131	130	0.003	ng/u1	0.02
46) Dimethylphthalate-d6	13.798	166	105	0.001	ng/u1	0.01
49) Acenaphthylene-d8	14.069	160	1074	0.012	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	54	0.003	ng/u1	0.00
60) Fluorene-d10	15.381	176	328	0.005	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	12	0.001	ng/u1	0.01
73) Anthracene-d10	17.280	188	13935	0.136	ng/u1	0.01
81) Pyrene-d10	19.645	212	1045	0.010	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.704	264	769	0.008	ng/u1	-0.02

Target Compounds	Qvalue

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

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