

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012725\
 Data File : BP023759.D
 Acq On : 28 Jan 2025 04:02
 Operator : RC/JU
 Sample : Q1174-21
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E2951

Quant Time: Jan 28 04:25:28 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012225.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 27 21:53:59 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	579653	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	2416254	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	1528663	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.216	188	3125764	20.000	ng/u1	0.00
79) Chrysene-d12	21.663	240	3146450	20.000	ng/u1	-0.01
88) Perylene-d12	25.080	264	3398037	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	64695	4.495	ng/uL	0.00
4) Pyridine-d5	3.711	84	182861	4.403	ng/u1	0.00
7) Phenol-d5	6.975	99	416936	8.501	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.122	67	909133	32.816	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	1198348	32.138	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	829266	21.501	ng/u1	0.02
21) Nitrobenzene-d5	8.934	128	663049	36.422	ng/u1	0.00
24) 2-Nitrophenol-d4	9.646	143	702419	42.086	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	1264207	36.036	ng/u1	0.02
31) 4-Chloroaniline-d4	10.699	131	1286783	22.936	ng/u1	0.00
46) Dimethylphthalate-d6	13.828	166	3991696	36.309	ng/u1	0.01
49) Acenaphthylene-d8	14.110	160	4553333	35.786	ng/u1	0.00
54) 4-Nitrophenol-d4	14.651	143	153919	6.968	ng/u1	0.06
60) Fluorene-d10	15.428	176	3533021	37.542	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.540	200	486380	31.598	ng/u1	0.01
73) Anthracene-d10	17.316	188	5697099	37.964	ng/u1	0.00
81) Pyrene-d10	19.692	212	6627146	41.392	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.845	264	6986201	41.075	ng/u1	0.00

Target Compounds	Qvalue

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

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