

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012725\
 Data File : BP023745.D
 Acq On : 27 Jan 2025 17:53
 Operator : RC/JU
 Sample : Q1174-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E2936

Quant Time: Jan 27 22:01:49 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	7.793	152	565692	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	2421148	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	1595777	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.216	188	3262566	20.000	ng/u1	0.00
79) Chrysene-d12	21.669	240	3297980	20.000	ng/u1	0.00
88) Perylene-d12	25.080	264	3557003	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	65008	4.629	ng/uL	0.00
4) Pyridine-d5	3.711	84	244675	6.037	ng/u1	0.00
7) Phenol-d5	6.975	99	384330	8.029	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.122	67	843575	31.201	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	1136732	31.238	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	765438	20.336	ng/u1	0.02
21) Nitrobenzene-d5	8.934	128	644101	35.309	ng/u1	0.00
24) 2-Nitrophenol-d4	9.658	143	699162	41.806	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.205	165	1245892	35.442	ng/u1	0.02
31) 4-Chloroaniline-d4	10.705	131	1493477	26.566	ng/u1	0.01
46) Dimethylphthalate-d6	13.828	166	4254480	37.071	ng/u1	0.01
49) Acenaphthylene-d8	14.116	160	4742129	35.703	ng/u1	0.01
54) 4-Nitrophenol-d4	14.646	143	166113	7.204	ng/u1	0.05
60) Fluorene-d10	15.428	176	3643008	37.083	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.540	200	574590	35.763	ng/u1	0.01
73) Anthracene-d10	17.322	188	6074688	38.783	ng/u1	0.00
81) Pyrene-d10	19.686	212	6842222	40.772	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.839	264	7182670	40.343	ng/u1	0.00

Target Compounds Qvalue

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

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