

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040125\
 Data File : BP024206.D
 Acq On : 01 Apr 2025 10:37
 Operator : RC/JU
 Sample : PB166622BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK622

Quant Time: Apr 01 12:37:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 01 12:35:32 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	353157	20.000	ng/u1	0.00
20) Naphthalene-d8	10.522	136	1527261	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.381	164	1000646	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	2042267	20.000	ng/u1	0.00
79) Chrysene-d12	21.639	240	2170251	20.000	ng/u1	0.00
88) Perylene-d12	25.015	264	2213165	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	61740	6.964	ng/uL	0.00
4) Pyridine-d5	3.675	84	850409	33.345	ng/u1	0.00
7) Phenol-d5	6.922	99	1044937	33.349	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.075	67	609194	35.304	ng/u1	0.00
11) 2-Chlorophenol-d4	7.281	132	832421	34.607	ng/u1	0.00
15) 4-Methylphenol-d8	8.446	113	821945	33.223	ng/u1	0.00
21) Nitrobenzene-d5	8.893	128	405095	32.849	ng/u1	0.00
24) 2-Nitrophenol-d4	9.604	143	426284	33.110	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.146	165	726919	31.400	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1131722	31.985	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2562074	35.334	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2755733	32.766	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	400148	26.955	ng/u1	0.00
60) Fluorene-d10	15.386	176	2104505	34.160	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	356908	27.680	ng/u1	0.00
73) Anthracene-d10	17.286	188	3437583	34.491	ng/u1	0.00
81) Pyrene-d10	19.669	212	4064644	37.383	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.780	264	4103865	35.767	ng/u1	-0.02

Target Compounds	Qvalue

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

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