

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100124\
 Data File : BP022130.D
 Acq On : 01 Oct 2024 11:06
 Operator : RC/JU
 Sample : PB163565BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK565

Quant Time: Oct 02 07:10:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	128962	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.451	136	461536	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.304	164	348762	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.098	188	817572	20.000	ng/u1	-0.02
79) Chrysene-d12	21.533	240	983899	20.000	ng/u1	0.00
88) Perylene-d12	24.815	264	1142448	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	15538	4.719	ng/uL	0.00
4) Pyridine-d5	3.628	84	209176	22.227	ng/u1	0.00
7) Phenol-d5	6.881	99	243942	23.124	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.034	67	152875	23.534	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.240	132	193681	25.376	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	196546	23.513	ng/u1	-0.01
21) Nitrobenzene-d5	8.834	128	91496	26.487	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.551	143	108831	27.786	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.081	165	206246	24.247	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.587	131	248256	24.461	ng/u1	-0.02
46) Dimethylphthalate-d6	13.704	166	683233	25.343	ng/u1	-0.02
49) Acenaphthylene-d8	13.992	160	713496	24.997	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.510	143	98651	21.393	ng/u1	-0.02
60) Fluorene-d10	15.304	176	579810	24.313	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.428	200	113562	22.373	ng/u1	-0.02
73) Anthracene-d10	17.198	188	976474	25.584	ng/u1	-0.02
81) Pyrene-d10	19.580	212	1279351	25.098	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.586	264	1485289	24.718	ng/u1	-0.02

Target Compounds Qvalue

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

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