

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091324\
 Data File : BP021905.D
 Acq On : 13 Sep 2024 12:26
 Operator : RC/JU
 Sample : PB163298TB 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB298

Quant Time: Sep 13 13:03:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 12 12:48:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	89273	20.000	ng/u1	0.03
20) Naphthalene-d8	10.534	136	497533	20.000	ng/u1	0.03
38) Acenaphthene-d10	14.369	164	360718	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	867384	20.000	ng/u1	0.00
79) Chrysene-d12	21.598	240	910739	20.000	ng/u1	0.00
88) Perylene-d12	24.921	264	988139	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	19233	6.969	ng/uL	0.01
4) Pyridine-d5	3.687	84	227744	28.028	ng/u1	0.02
7) Phenol-d5	6.940	99	209157	27.420	ng/u1	0.03
9) Bis-(2-Chloroethyl)eth...	7.099	67	131942	29.780	ng/u1	0.03
11) 2-Chlorophenol-d4	7.298	132	171372	32.274	ng/u1	0.02
15) 4-Methylphenol-d8	8.457	113	173475	28.484	ng/u1	0.03
21) Nitrobenzene-d5	8.904	128	87441	28.219	ng/u1	0.03
24) 2-Nitrophenol-d4	9.616	143	124503	35.345	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.151	165	247255	31.697	ng/u1	0.02
31) 4-Chloroaniline-d4	10.657	131	325788	28.657	ng/u1	0.02
46) Dimethylphthalate-d6	13.769	166	920126	33.953	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	930858	31.313	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	136868	26.716	ng/u1	0.00
60) Fluorene-d10	15.375	176	782268	30.369	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	122192	24.624	ng/u1	0.00
73) Anthracene-d10	17.263	188	1334288	33.116	ng/u1	0.00
81) Pyrene-d10	19.633	212	1678244	33.306	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.692	264	1679603	32.661	ng/u1	0.00

Target Compounds	Qvalue

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

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