

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022074.D
 Acq On : 27 Sep 2024 10:52
 Operator : RC/JU
 Sample : PB163390BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK390

Quant Time: Sep 28 03:06:59 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	7.681	152	114759	20.000	ng/u1	-0.03
20) Naphthalene-d8	10.440	136	432198	20.000	ng/u1	-0.03
38) Acenaphthene-d10	14.304	164	319893	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.116	188	794718	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	904246	20.000	ng/u1	0.00
88) Perylene-d12	24.833	264	991967	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	15979	5.454	ng/uL	-0.02
4) Pyridine-d5	3.617	84	213443	25.487	ng/u1	-0.02
7) Phenol-d5	6.864	99	251922	26.836	ng/u1	-0.03
9) Bis-(2-Chloroethyl)eth...	7.016	67	158296	27.385	ng/u1	-0.03
11) 2-Chlorophenol-d4	7.222	132	188200	27.709	ng/u1	-0.02
15) 4-Methylphenol-d8	8.381	113	196487	26.415	ng/u1	-0.02
21) Nitrobenzene-d5	8.816	128	92956	28.736	ng/u1	-0.03
24) 2-Nitrophenol-d4	9.534	143	106496	29.035	ng/u1	-0.03
28) 2,4-Dichlorophenol-d3	10.069	165	217649	27.325	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.581	131	235517	24.781	ng/u1	-0.02
46) Dimethylphthalate-d6	13.710	166	729541	29.502	ng/u1	-0.01
49) Acenaphthylene-d8	13.998	160	742177	28.348	ng/u1	0.00
54) 4-Nitrophenol-d4	14.510	143	101552	24.009	ng/u1	-0.02
60) Fluorene-d10	15.316	176	624078	28.531	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.440	200	102752	20.825	ng/u1	0.00
73) Anthracene-d10	17.210	188	1043156	28.117	ng/u1	0.00
81) Pyrene-d10	19.598	212	1422996	30.375	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.616	264	1526387	29.256	ng/u1	0.00

Target Compounds Qvalue

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

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