

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
 Data File : BP021833.D
 Acq On : 05 Sep 2024 03:04
 Operator : RC/JU
 Sample : PB163099BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK099

Quant Time: Sep 05 06:18:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 05 05:03:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	293408	20.000	ng/u1	0.00
20) Naphthalene-d8	10.557	136	1182120	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.410	164	690290	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.210	188	1595859	20.000	ng/u1	0.00
79) Chrysene-d12	21.657	240	1604810	20.000	ng/u1	-0.02
88) Perylene-d12	25.068	264	1729903	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	52617	5.954	ng/uL	0.00
4) Pyridine-d5	3.681	84	689215	26.962	ng/u1	0.00
7) Phenol-d5	6.946	99	855735	27.329	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.105	67	477387	28.015	ng/u1	0.01
11) 2-Chlorophenol-d4	7.305	132	570279	28.210	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	644837	26.650	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	285455	30.014	ng/u1	0.00
24) 2-Nitrophenol-d4	9.640	143	293519	30.404	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.175	165	524383	27.564	ng/u1	0.00
31) 4-Chloroaniline-d4	10.687	131	739791	26.729	ng/u1	0.00
46) Dimethylphthalate-d6	13.810	166	1842629	34.886	ng/u1	0.00
49) Acenaphthylene-d8	14.098	160	1946458	32.823	ng/u1	0.00
54) 4-Nitrophenol-d4	14.622	143	287798	28.640	ng/u1	0.01
60) Fluorene-d10	15.416	176	1391341	31.332	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.534	200	217409	24.769	ng/u1	0.00
73) Anthracene-d10	17.316	188	2293246	30.279	ng/u1	0.00
81) Pyrene-d10	19.680	212	2769265	30.187	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.821	264	2791745	30.265	ng/u1	-0.04

Target Compounds Qvalue

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

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