

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017063.D
 Acq On : 09 Sep 2023 22:05
 Operator : MA/JU
 Sample : PB155238BL
 Misc :
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK238

Quant Time: Sep 10 00:56:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 01:41:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	560877	20.000	ng/u1	0.00
20) Naphthalene-d8	10.793	136	2304485	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.628	164	1278817	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	2306887	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	1444236	20.000	ng/u1	0.00
88) Perylene-d12	24.015	264	1480885	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	80620	5.437	ng/uL	0.00
4) Pyridine-d5	3.764	84	617950	14.585	ng/u1	0.00
7) Phenol-d5	7.116	99	1371133	26.758	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	930391	29.028	ng/u1	0.00
11) 2-Chlorophenol-d4	7.487	132	1057266	28.509	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	1056847	27.073	ng/u1	-0.01
21) Nitrobenzene-d5	9.169	128	508400	28.437	ng/u1	0.00
24) 2-Nitrophenol-d4	9.881	143	530904	27.394	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	920598	25.729	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	466399	9.035	ng/u1	0.00
46) Dimethylphthalate-d6	14.040	166	2688748	28.246	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	3123083	28.429	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	397636	21.385	ng/u1	0.00
60) Fluorene-d10	15.622	176	2221242	27.254	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	291410	20.991	ng/u1	-0.01
73) Anthracene-d10	17.492	188	3144784	29.684	ng/u1	0.00
81) Pyrene-d10	19.727	212	2975641	33.826	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	2300186	30.617	ng/u1	0.00

Target Compounds Qvalue

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

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