

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017478.D
 Acq On : 02 Oct 2023 15:41
 Operator : MA/JU
 Sample : PB155899BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK899

Quant Time: Oct 02 23:21:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	94545	20.000	ng/u1	0.00
20) Naphthalene-d8	10.787	136	376541	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.639	164	220905	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.427	188	494032	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	448402	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	482541	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	16447	7.145	ng/uL	0.00
4) Pyridine-d5	3.728	84	222628	34.597	ng/u1	0.00
7) Phenol-d5	7.104	99	271252	34.933	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	168435	35.943	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	223266	36.253	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	211981	35.114	ng/u1	0.00
21) Nitrobenzene-d5	9.157	128	104894	38.533	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	109657	37.099	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	206245	36.933	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	290600	35.924	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	625420	39.521	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	696678	38.278	ng/u1	0.00
54) 4-Nitrophenol-d4	14.880	143	53253	19.316	ng/u1	0.00
60) Fluorene-d10	15.645	176	541244	39.728	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	54966	19.740	ng/u1	0.00
73) Anthracene-d10	17.527	188	875781	39.971	ng/u1	0.00
81) Pyrene-d10	19.780	212	1054203	43.475	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.956	264	911605	39.238	ng/u1	0.00

Target Compounds Qvalue

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

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