

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
 Data File : BP017094.D
 Acq On : 11 Sep 2023 08:43
 Operator : MA/JU
 Sample : PB155284BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK284

Quant Time: Sep 11 09:18:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 06:13:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	373976	20.000	ng/u1	0.00
20) Naphthalene-d8	10.787	136	1730278	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1147182	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	2480478	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	2054595	20.000	ng/u1	0.00
88) Perylene-d12	24.015	264	2113330	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	59219	6.044	ng/uL	0.00
4) Pyridine-d5	3.752	84	867956	29.376	ng/u1	0.00
7) Phenol-d5	7.111	99	993845	29.235	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	694497	30.461	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	751604	30.591	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	776577	28.878	ng/u1	0.00
21) Nitrobenzene-d5	9.157	128	369481	29.994	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	396278	31.169	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	683063	28.334	ng/u1	0.00
31) 4-Chloroaniline-d4	10.946	131	1086354	28.102	ng/u1	0.00
46) Dimethylphthalate-d6	14.034	166	2455642	30.083	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	2590769	27.873	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	424928	25.232	ng/u1	0.00
60) Fluorene-d10	15.616	176	2006124	28.961	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	315697	22.553	ng/u1	0.00
73) Anthracene-d10	17.486	188	3208528	30.178	ng/u1	0.00
81) Pyrene-d10	19.727	212	3620051	33.458	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	3104912	30.604	ng/u1	0.00

Target Compounds Qvalue

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

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